

AP-111

Investigation Report  
(SWMU-10 Sludge  
Pits)  
(3 of 3)

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605875

07-Jul-16

Client: Western Refining Company

Project: SWMU 10

|                            |           |      |                          |             |      |   |           |              |          |      |
|----------------------------|-----------|------|--------------------------|-------------|------|---|-----------|--------------|----------|------|
| Sample ID                  | Ics-25469 |      | SampType: LCS            |             |      | TestCode: EPA Method 8270C: Semivolatiles |           |              |          |      |
| Client ID:                 | LCSS      |      | Batch ID: 25469          |             |      | RunNo: 34461                              |           |              |          |      |
| Prep Date:                 | 5/24/2016 |      | Analysis Date: 5/25/2016 |             |      | SeqNo: 1063512                            |           | Units: mg/Kg |          |      |
| Analyte                    | Result    | PQL  | SPK value                | SPK Ref Val | %REC | LowLimit                                  | HighLimit | %RPD         | RPDLimit | Qual |
| Acenaphthene               | 1.2       | 0.20 | 1.670                    | 0           | 70.0 | 45.8                                      | 99.8      |              |          |      |
| 4-Chloro-3-methylphenol    | 2.2       | 0.50 | 3.330                    | 0           | 66.5 | 51.5                                      | 103       |              |          |      |
| 2-Chlorophenol             | 1.8       | 0.20 | 3.330                    | 0           | 54.1 | 46.5                                      | 105       |              |          |      |
| 1,4-Dichlorobenzene        | 0.86      | 0.20 | 1.670                    | 0           | 51.3 | 45.5                                      | 103       |              |          |      |
| 2,4-Dinitrotoluene         | 1.1       | 0.50 | 1.670                    | 0           | 66.9 | 36  | 87.2      |              |          |      |
| N-Nitrosodi-n-propylamine  | 0.80      | 0.20 | 1.670                    | 0           | 48.1 | 47.3                                      | 104       |              |          |      |
| 4-Nitrophenol              | 2.2       | 0.25 | 3.330                    | 0           | 65.7 | 47.3                                      | 95.3      |              |          |      |
| Pentachlorophenol          | 1.9       | 0.40 | 3.330                    | 0           | 56.6 | 38.7                                      | 89.3      |              |          |      |
| Phenol                     | 1.8       | 0.20 | 3.330                    | 0           | 54.5 | 47.8                                      | 106       |              |          |      |
| Pyrene                     | 1.2       | 0.20 | 1.670                    | 0           | 72.6 | 33.4                                      | 105       |              |          |      |
| 1,2,4-Trichlorobenzene     | 1.0       | 0.20 | 1.670                    | 0           | 61.4 | 50.4                                      | 115       |              |          |      |
| Surr: 2-Fluorophenol       | 1.6       |      | 3.330                    |             | 47.4 | 28.3                                      | 102       |              |          |      |
| Surr: Phenol-d5            | 1.9       |      | 3.330                    |             | 56.0 | 35.7                                      | 103       |              |          |      |
| Surr: 2,4,6-Tribromophenol | 2.2       |      | 3.330                    |             | 66.3 | 35.2                                      | 108       |              |          |      |
| Surr: Nitrobenzene-d5      | 1.0       |      | 1.670                    |             | 59.7 | 24  | 118       |              |          |      |
| Surr: 2-Fluorobiphenyl     | 1.2       |      | 1.670                    |             | 70.2 | 35.4                                      | 111       |              |          |      |
| Surr: 4-Terphenyl-d14      | 0.74      |      | 1.670                    |             | 44.5 | 15  | 91.7      |              |          |      |

### Qualifiers:

|    |   |    |   |
|----|---|----|---|
| *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
| D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
| H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
| ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
| R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
| S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

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|------------|-----------|-------|----------------|-------------|------|-----------|--------------------------|------|----------|-------|--|
| Sample ID  | MB-25448  |       | SampType:      | MBLK        |      | TestCode: | EPA Method 7471: Mercury |      |          |       |  |
| Client ID: | PBS       |       | Batch ID:      | 25448       |      | RunNo:    | 34452                    |      |          |       |  |
| Prep Date: | 5/23/2016 |       | Analysis Date: | 5/24/2016   |      | SeqNo:    | 1062465                  |      | Units:   | mg/Kg |  |
| Analyte    | Result    | PQL   | SPK value      | SPK Ref Val | %REC | LowLimit  | HighLimit                | %RPD | RPDLimit | Qual  |  |
| Mercurv    | 0.00083   | 0.033 |                |             |      |           |                          |      |          | J     |  |

|            |           |       |           |                |           |          |           |                          |          |       |  |
|------------|-----------|-------|-----------|----------------|-----------|----------|-----------|--------------------------|----------|-------|--|
| Sample ID  | LCS-25448 |       |           | SampType:      | LCS       |          | TestCode: | EPA Method 7471: Mercury |          |       |  |
| Client ID: | LCSS      |       |           | Batch ID:      | 25448     |          | RunNo:    | 34452                    |          |       |  |
| Prep Date: | 5/23/2016 |       |           | Analysis Date: | 5/24/2016 |          | SeqNo:    | 1062466                  | Units:   | mg/Kg |  |
| Analyte    | Result    | PQL   | SPK value | SPK Ref Val    | %REC      | LowLimit | HighLimit | %RPD                     | RPDLimit | Qual  |  |
| Mercury    | 0.18      | 0.033 | 0.1667    | 0              | 110       | 80       | 120       |                          |          |       |  |

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D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
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P Sample pH Not In Range  
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W Sample container temperature is out of limit as specified

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|------------|------------------|------|----------------|------------------|------|-----------|--------------------------------------|------|---------------------|------|
| Sample ID  | <b>MB-25447</b>  |      | SampType:      | <b>MBLK</b>      |      | TestCode: | <b>EPA Method 6010B: Soil Metals</b> |      |                     |      |
| Client ID: | <b>PBS</b>       |      | Batch ID:      | <b>25447</b>     |      | RunNo:    | <b>34439</b>                         |      |                     |      |
| Prep Date: | <b>5/23/2016</b> |      | Analysis Date: | <b>5/24/2016</b> |      | SeqNo:    | <b>1061954</b>                       |      | Units: <b>mg/Kg</b> |      |
| Analyte    | Result           | PQL  | SPK value      | SPK Ref Val      | %REC | LowLimit  | HighLimit                            | %RPD | RPDLimit            | Qual |
| Antimony   | ND               | 2.5  |                |                  |      |           |                                      |      |                     |      |
| Arsenic    | ND               | 2.5  |                |                  |      |           |                                      |      |                     |      |
| Beryllium  | ND               | 0.15 |                |                  |      |           |                                      |      |                     |      |
| Cadmium    | ND               | 0.10 |                |                  |      |           |                                      |      |                     |      |
| Chromium   | ND               | 0.30 |                |                  |      |           |                                      |      |                     |      |
| Cobalt     | ND               | 0.30 |                |                  |      |           |                                      |      |                     |      |
| Lead       | ND               | 0.25 |                |                  |      |           |                                      |      |                     |      |
| Manganese  | 0.047            | 0.10 |                |                  |      |           |                                      |      |                     | J    |
| Nickel     | ND               | 0.50 |                |                  |      |           |                                      |      |                     |      |
| Selenium   | 1.3              | 2.5  |                |                  |      |           |                                      |      |                     | J    |
| Silver     | ND               | 0.25 |                |                  |      |           |                                      |      |                     |      |
| Vanadium   | ND               | 2.5  |                |                  |      |           |                                      |      |                     |      |
| Zinc       | ND               | 2.5  |                |                  |      |           |                                      |      |                     |      |

|            |                  |      |                |                  |      |           |                                      |      |                     |      |
|------------|------------------|------|----------------|------------------|------|-----------|--------------------------------------|------|---------------------|------|
| Sample ID  | <b>LCS-25447</b> |      | SampType:      | <b>LCS</b>       |      | TestCode: | <b>EPA Method 6010B: Soil Metals</b> |      |                     |      |
| Client ID: | <b>LCSS</b>      |      | Batch ID:      | <b>25447</b>     |      | RunNo:    | <b>34439</b>                         |      |                     |      |
| Prep Date: | <b>5/23/2016</b> |      | Analysis Date: | <b>5/24/2016</b> |      | SeqNo:    | <b>1061955</b>                       |      | Units: <b>mg/Kg</b> |      |
| Analyte    | Result           | PQL  | SPK value      | SPK Ref Val      | %REC | LowLimit  | HighLimit                            | %RPD | RPDLimit            | Qual |
| Antimony   | 23               | 2.5  | 25.00          | 0                | 90.0 | 80        | 120                                  |      |                     |      |
| Arsenic    | 24               | 2.5  | 25.00          | 0                | 95.3 | 80        | 120                                  |      |                     |      |
| Beryllium  | 24               | 0.15 | 25.00          | 0                | 97.0 | 80        | 120                                  |      |                     |      |
| Cadmium    | 24               | 0.10 | 25.00          | 0                | 94.4 | 80        | 120                                  |      |                     |      |
| Chromium   | 23               | 0.30 | 25.00          | 0                | 93.1 | 80        | 120                                  |      |                     |      |
| Cobalt     | 23               | 0.30 | 25.00          | 0                | 90.9 | 80        | 120                                  |      |                     |      |
| Lead       | 23               | 0.25 | 25.00          | 0                | 91.3 | 80        | 120                                  |      |                     |      |
| Manganese  | 24               | 0.10 | 25.00          | 0                | 94.1 | 80        | 120                                  |      |                     |      |
| Nickel     | 23               | 0.50 | 25.00          | 0                | 92.2 | 80        | 120                                  |      |                     |      |
| Selenium   | 24               | 2.5  | 25.00          | 0                | 97.6 | 80        | 120                                  |      |                     |      |
| Silver     | 4.7              | 0.25 | 5.000          | 0                | 93.8 | 80        | 120                                  |      |                     |      |
| Vanadium   | 24               | 2.5  | 25.00          | 0                | 95.9 | 80        | 120                                  |      |                     |      |
| Zinc       | 24               | 2.5  | 25.00          | 0                | 94.2 | 80        | 120                                  |      |                     |      |

|            |                  |      |                |                 |      |           |                                      |      |                     |      |
|------------|------------------|------|----------------|-----------------|------|-----------|--------------------------------------|------|---------------------|------|
| Sample ID  | <b>MB-25586</b>  |      | SampType:      | <b>MBLK</b>     |      | TestCode: | <b>EPA Method 6010B: Soil Metals</b> |      |                     |      |
| Client ID: | <b>PBS</b>       |      | Batch ID:      | <b>25586</b>    |      | RunNo:    | <b>34597</b>                         |      |                     |      |
| Prep Date: | <b>5/31/2016</b> |      | Analysis Date: | <b>6/1/2016</b> |      | SeqNo:    | <b>1067058</b>                       |      | Units: <b>mg/Kg</b> |      |
| Analyte    | Result           | PQL  | SPK value      | SPK Ref Val     | %REC | LowLimit  | HighLimit                            | %RPD | RPDLimit            | Qual |
| Barium     | ND               | 0.10 |                |                 |      |           |                                      |      |                     |      |

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|------------|-----------|------|-------------------------|-------------|---|----------|--------------|------|----------|------|
| Sample ID  | LCS-25586 |      | SampType: LCS           |             | TestCode: EPA Method 6010B: Soil Metals |          |              |      |          |      |
| Client ID: | LCSS      |      | Batch ID: 25586         |             | RunNo: 34597                            |          |              |      |          |      |
| Prep Date: | 5/31/2016 |      | Analysis Date: 6/1/2016 |             | SeqNo: 1067059                          |          | Units: mg/Kg |      |          |      |
| Analyte    | Result    | PQL  | SPK value               | SPK Ref Val | %REC                                    | LowLimit | HighLimit    | %RPD | RPDLimit | Qual |
| Barium     | 22        | 0.10 | 25.00                   | 0           | 88.8                                    | 80       | 120          |      |          |      |

|            |           |     |                          |             |   |          |              |      |          |      |
|------------|-----------|-----|--------------------------|-------------|---|----------|--------------|------|----------|------|
| Sample ID  | MB-25586  |     | SampType: MBLK           |             | TestCode: EPA Method 6010B: Soil Metals |          |              |      |          |      |
| Client ID: | PBS       |     | Batch ID: 25586          |             | RunNo: 34951                            |          |              |      |          |      |
| Prep Date: | 5/31/2016 |     | Analysis Date: 6/16/2016 |             | SeqNo: 1079711                          |          | Units: mg/Kg |      |          |      |
| Analyte    | Result    | PQL | SPK value                | SPK Ref Val | %REC                                    | LowLimit | HighLimit    | %RPD | RPDLimit | Qual |
| Iron       | 2.0       | 2.5 |                          |             |   |          |              |      |          | J    |

|            |           |     |                          |             |   |          |              |      |          |      |
|------------|-----------|-----|--------------------------|-------------|---|----------|--------------|------|----------|------|
| Sample ID  | LCS-25586 |     | SampType: LCS            |             | TestCode: EPA Method 6010B: Soil Metals |          |              |      |          |      |
| Client ID: | LCSS      |     | Batch ID: 25586          |             | RunNo: 34951                            |          |              |      |          |      |
| Prep Date: | 5/31/2016 |     | Analysis Date: 6/16/2016 |             | SeqNo: 1079712                          |          | Units: mg/Kg |      |          |      |
| Analyte    | Result    | PQL | SPK value                | SPK Ref Val | %REC                                    | LowLimit | HighLimit    | %RPD | RPDLimit | Qual |
| Iron       | 25        | 2.5 | 25.00                    | 0           | 99.4                                    | 80       | 120          |      |          |      |

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|-------------------------------|--------|----------------|-----------|-------------|--------------------------------------|----------|-----------|------|----------|------|
| Sample ID                     | rb     | SampType:      | MBLK      | TestCode:   | EPA Method 8015D Mod: Gasoline Range |          |           |      |          |      |
| Client ID:                    | PBS    | Batch ID:      | C34413    | RunNo:      | 34413                                |          |           |      |          |      |
| Prep Date:                    |        | Analysis Date: | 5/23/2016 | SeqNo:      | 1061608                              | Units:   | mg/Kg     |      |          |      |
| Analyte                       | Result | PQL            | SPK value | SPK Ref Val | %REC                                 | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO) | ND     | 5.0            |           |             |                                      |          |           |      |          |      |
| Surr: BFB                     | 500    |                | 500.0     |             | 100                                  | 70       | 130       |      |          |      |

|                               |               |                |           |             |                                      |          |           |      |          |      |
|-------------------------------|---------------|----------------|-----------|-------------|--------------------------------------|----------|-----------|------|----------|------|
| Sample ID                     | 2.5ug gro icv | SampType:      | LCS       | TestCode:   | EPA Method 8015D Mod: Gasoline Range |          |           |      |          |      |
| Client ID:                    | LCSS          | Batch ID:      | C34413    | RunNo:      | 34413                                |          |           |      |          |      |
| Prep Date:                    |               | Analysis Date: | 5/23/2016 | SeqNo:      | 1061609                              | Units:   | mg/Kg     |      |          |      |
| Analyte                       | Result        | PQL            | SPK value | SPK Ref Val | %REC                                 | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO) | 26            | 5.0            | 25.00     | 0           | 102                                  | 62.9     | 123       |      |          |      |
| Surr: BFB                     | 510           |                | 500.0     |             | 102                                  | 70       | 130       |      |          |      |

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## Sample Log-In Check List

Client Name: Western Refining Gallup

Work Order Number: 1605375

RcptNo: 1

Received by/date:

SA 05/18/16

Logged By: Lindsay Mangin

5/18/2016 4:00:00 PM

*[Signature]*

Completed By: Lindsay Mangin

5/19/2016 7:00:19 AM

*[Signature]*

Reviewed By:

*[Signature]* 05/19/16

### Chain of Custody

1. Custody seals intact on sample bottles? Yes ☐ No ☐ Not Present ☒
2. Is Chain of Custody complete? Yes ☒ No ☐ Not Present ☐
3. How was the sample delivered? Courier

### Log In

4. Was an attempt made to cool the samples? Yes ☒ No ☐ NA ☐
5. Were all samples received at a temperature of  $>0^{\circ}\text{C}$  to  $6.0^{\circ}\text{C}$ ? Yes ☒ No ☐ NA ☐
6. Sample(s) in proper container(s)? Yes ☒ No ☐
7. Sufficient sample volume for indicated test(s)? Yes ☒ No ☐
8. Are samples (except VOA and ONG) properly preserved? Yes ☒ No ☐
9. Was preservative added to bottles? Yes ☐ No ☒ NA ☐
10. VOA vials have zero headspace? Yes ☐ No ☐ No VOA Vials ☒
11. Were any sample containers received broken? Yes ☐ No ☒
12. Does paperwork match bottle labels?  
(Note discrepancies on chain of custody) Yes ☒ No ☐
13. Are matrices correctly identified on Chain of Custody? Yes ☒ No ☐
14. Is it clear what analyses were requested? Yes ☒ No ☐
15. Were all holding times able to be met?  
(If no, notify customer for authorization.) Yes ☒ No ☐

# of preserved  
bottles checked  
for pH:

( $<2$  or  $>12$  unless noted)

Adjusted? \_\_\_\_\_

Checked by: \_\_\_\_\_

### Special Handling (if applicable)

16. Was client notified of all discrepancies with this order? Yes ☐ No ☐ NA ☒

Person Notified: \_\_\_\_\_

Date: \_\_\_\_\_

By Whom: \_\_\_\_\_

Via: ☐ eMail ☐ Phone ☐ Fax ☐ In Person

Regarding: \_\_\_\_\_

Client Instructions: \_\_\_\_\_

17. Additional remarks:

### 18. Cooler Information

| Cooler No | Temp $^{\circ}\text{C}$ | Condition | Seal Intact | Seal No | Seal Date | Signed By |
|-----------|-------------------------|-----------|-------------|---------|-----------|-----------|
| 1         | 1.0                     | Good      | Yes         |         |           |           |

# Chain-of-Custody Record

Client: WESTERN REFINING SW, INC.  
GALLUP REFINERY  
 Mailing Address: 92 GIANT CROSSING RD  
GALLUP, NM 87301  
 Phone #: 505-722-0217  
 Mail or Fax#: ED.RIEGE@WNR.COM

A/QC Package:  
☒ Standard ☐ Level 4 (Full Validation)  
 Accreditation  
☒ NELAP ☐ Other

EDD (Type) EXCEL

| Date  | Time | Matrix | Sample Request ID   | Container Type and # | Preservative Type | HEAL No. |
|-------|------|--------|---------------------|----------------------|-------------------|----------|
| 17-16 | 1325 | SOIL   | SWMU 10-20 (8-10')  | JARS-3               | NEAT              | 11055875 |
|       |      |        |                     | VIALS-2              | MEOH              | -001     |
|       |      |        |                     | VIALS-2              | SOBI              | -001     |
| 1610  |      |        | SWMU 10-20 (10-12') | JARS-3               | NEAT              | -002     |
|       |      |        |                     | VIALS-2              | MEOH              | -002     |
|       |      |        |                     | VIALS-2              | SOBI              | -002     |
| 1625  |      |        | SWMU 10-20 (16-18') | JARS-3               | NEAT              | -003     |
|       |      |        |                     | VIALS-2              | MEOH              | -003     |
|       |      |        |                     | VIALS-2              | SOBI              | -003     |
| 1635  |      |        | SWMU 10-20 (20-22') | JARS-3               | NEAT              | -004     |
|       |      |        |                     | VIALS-2              | MEOH              | -004     |
|       |      |        |                     | VIALS-2              | SOBI              | -004     |

Relinquished by: AF Date: 5/18/16 Time: 800  
 Relinquished by: AF Date: 5/18/16 Time: 1600

Turn-Around Time:

☒ Standard ☐ Rush  
 Project Name:

SWMU 10  
 Project #:

Project Manager:

ED RIEGE  
 Sampler: TRACY PAYNE  
 On Ice: ☒ Yes ☐ No

Sample Temperature: 1.0

## HALL ENVIRONMENTAL ANALYSIS LABORATORY

www.hallenvironmental.com  
 4901 Hawkins NE - Albuquerque, NM 87109  
 Tel. 505-345-3975 Fax 505-345-4107

### Analysis Request

| BTEX + MTBE + TMBs (8021) | BTEX + MTBE + TPH (Gas only) | TPH 8015B (GRO / DRO / MRO) | TPH (Method 418.1) | EDB (Method 504.1) | PAH's (8310 or 8270 SIMS) | RORA 8 Metals | Anions (F, Cl, NO <sub>3</sub> , NO <sub>2</sub> , PO <sub>4</sub> , SO <sub>4</sub> ) | 8081 Pesticides / 8082 PCB's | 8260B (VOA) | 8270 (Semi-VOA) | METALS & CYANIDE | (SEE ATTACHED LIST) | Air Bubbles (Y or N) |
|---------------------------|------------------------------|-----------------------------|--------------------|--------------------|---------------------------|---------------|--|------------------------------|-------------|-----------------|------------------|---------------------|----------------------|
|                           |                              | ✓                           |                    |                    |                           |               |  |                              | ✓           | ✓               | ✓                |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              | ✓                           |                    |                    |                           |               |  |                              | ✓           | ✓               | ✓                |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              | ✓                           |                    |                    |                           |               |  |                              | ✓           | ✓               | ✓                |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              | ✓                           |                    |                    |                           |               |  |                              | ✓           | ✓               | ✓                |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |
|                           |                              |                             |                    |                    |                           |               |  |                              | ✓           |                 |                  |                     |                      |

Remarks: PLEASE INCLUDE J FLAGS AND MDL



**WESTERN REFINING SOUTHWEST, INC.**  
**GALLUP REFINERY - SWMU 10**  
**METALS AND CYANIDE ANALYSES FOR SOIL SAMPLES**

| <b>Analyte</b> | <b>Analytical Method</b>      |
|----------------|-------------------------------|
| Antimony       | SW-846 method 6010/6020       |
| Arsenic        | SW-846 method 6010/6020       |
| Barium         | SW-846 method 6010/6020       |
| Beryllium      | SW-846 method 6010/6020       |
| Cadmium        | SW-846 method 6010/6020       |
| Chromium       | SW-846 method 6010/6020       |
| Chromium VI    | SW-846 Method 3060A           |
| Cobalt         | SW-846 method 6010/6020       |
| Cyanide        | SW-846 method 335.4/335.2 mod |
| Lead           | SW-846 method 6010/6020       |
| Mercury        | SW-846 method 7470/7471       |
| Nickel         | SW-846 method 6010/6020       |
| Selenium       | SW-846 method 6010/6020       |
| Silver         | SW-846 method 6010/6020       |
| Vanadium       | SW-846 method 6010/6020       |
| Zinc           | SW-846 method 6010/6020       |
| Iron           | SW-846 method 6010/6020       |
| Manganese      | SW-846 method 6010/6020       |



Hall Environmental Analysis Laboratory  
4901 Hawkins NE  
Albuquerque, NM 87109  
TEL: 505-345-3975 FAX: 505-345-4107  
Website: [www.hallenvironmental.com](http://www.hallenvironmental.com)

June 28, 2016

Ed Riege  
Western Refining Company  
Rt. 3 Box 7  
Gallup, NM 87301  
TEL: (505) 722-0231  
FAX

RE: SWMU 10

OrderNo.: 1605943

Dear Ed Riege:

Hall Environmental Analysis Laboratory received 7 sample(s) on 5/19/2016 for the analyses presented in the following report.

These were analyzed according to EPA procedures or equivalent. To access our accredited tests please go to [www.hallenvironmental.com](http://www.hallenvironmental.com) or the state specific web sites. In order to properly interpret your results it is imperative that you review this report in its entirety. See the sample checklist and/or the Chain of Custody for information regarding the sample receipt temperature and preservation. Data qualifiers or a narrative will be provided if the sample analysis or analytical quality control parameters require a flag. When necessary, data qualifiers are provided on both the sample analysis report and the QC summary report, both sections should be reviewed. All samples are reported, as received, unless otherwise indicated. Lab measurement of analytes considered field parameters that require analysis within 15 minutes of sampling such as pH and residual chlorine are qualified as being analyzed outside of the recommended holding time.

Please don't hesitate to contact HEAL for any additional information or clarifications.

ADHS Cert #AZ0682 -- NMED-DWB Cert #NM9425 -- NMED-Micro Cert #NM0190

Sincerely,

A handwritten signature in black ink, appearing to read 'Andy Freeman'.

Andy Freeman  
Laboratory Manager  
4901 Hawkins NE  
Albuquerque, NM 87109

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-21 (2-2.5')

**Project:** SWMU 10

**Collection Date:** 5/18/2016 9:05:00 AM

**Lab ID:** 1605943-001

**Matrix:** SOIL

**Received Date:** 5/19/2016 4:30:00 PM

| Analyses   | Result | MDL     | PQL    | Qual | Units    | DF  | Date Analyzed         | Batch ID |
|--|--------|---------|--------|------|----------|-----|-----------------------|----------|
| <b>EPA METHOD 8015M/D: DIESEL RANGE ORGANICS</b> |        |         |        |      |          |     | Analyst: <b>KJH</b>   |          |
| Diesel Range Organics (DRO)                      | 32     | 3.7     | 10     |      | mg/Kg-dr | 1   | 5/24/2016 2:54:59 PM  | 25449    |
| Motor Oil Range Organics (MRO)                   | ND     | 52      | 52     |      | mg/Kg-dr | 1   | 5/24/2016 2:54:59 PM  | 25449    |
| Surr: DNOP                                       | 110    | 0       | 70-130 |      | %Rec     | 1   | 5/24/2016 2:54:59 PM  | 25449    |
| <b>PERCENT MOISTURE</b>                          |        |         |        |      |          |     | Analyst: <b>BCN</b>   |          |
| Percent Moisture                                 | 7.7    | 1.0     | 1.0    |      | wt%      | 1   | 5/23/2016             | R34494   |
| <b>EPA METHOD 7471: MERCURY</b>                  |        |         |        |      |          |     | Analyst: <b>pmf</b>   |          |
| Mercury  | 0.056  | 0.00062 | 0.036  |      | mg/Kg-dr | 1   | 5/26/2016 3:40:58 PM  | 25509    |
| <b>EPA METHOD 6010B: SOIL METALS</b>             |        |         |        |      |          |     | Analyst: <b>MED</b>   |          |
| Antimony   | ND     | 1.3     | 2.7    |      | mg/Kg-dr | 1   | 5/31/2016 9:51:45 AM  | 25508    |
| Arsenic  | 3.3    | 0.79    | 2.7    |      | mg/Kg-dr | 1   | 5/26/2016 4:04:00 PM  | 25508    |
| Barium   | 570    | 0.26    | 0.54   |      | mg/Kg-dr | 5   | 5/31/2016 10:06:38 AM | 25508    |
| Beryllium  | 0.72   | 0.0050  | 0.16   |      | mg/Kg-dr | 1   | 5/26/2016 4:04:00 PM  | 25508    |
| Cadmium  | ND     | 0.069   | 0.11   |      | mg/Kg-dr | 1   | 5/26/2016 4:04:00 PM  | 25508    |
| Chromium   | 9.2    | 0.14    | 0.33   |      | mg/Kg-dr | 1   | 5/26/2016 4:04:00 PM  | 25508    |
| Cobalt   | 3.7    | 0.041   | 0.33   |      | mg/Kg-dr | 1   | 5/26/2016 4:04:00 PM  | 25508    |
| Iron   | 9900   | 110     | 270    |      | mg/Kg-dr | 100 | 5/31/2016 10:08:25 AM | 25508    |
| Lead   | 2.1    | 0.19    | 0.27   |      | mg/Kg-dr | 1   | 5/31/2016 9:51:45 AM  | 25508    |
| Manganese  | 440    | 0.24    | 0.54   |      | mg/Kg-dr | 5   | 5/31/2016 10:06:38 AM | 25508    |
| Nickel   | 6.0    | 0.12    | 0.54   |      | mg/Kg-dr | 1   | 5/31/2016 9:51:45 AM  | 25508    |
| Selenium   | ND     | 1.2     | 2.7    |      | mg/Kg-dr | 1   | 5/31/2016 9:51:45 AM  | 25508    |
| Silver   | ND     | 0.034   | 0.27   |      | mg/Kg-dr | 1   | 5/26/2016 4:04:00 PM  | 25508    |
| Vanadium   | 16     | 0.038   | 2.7    |      | mg/Kg-dr | 1   | 5/26/2016 4:04:00 PM  | 25508    |
| Zinc   | 16     | 0.62    | 2.7    |      | mg/Kg-dr | 1   | 5/26/2016 4:04:00 PM  | 25508    |
| <b>EPA METHOD 8270C: SEMIVOLATILES</b>           |        |         |        |      |          |     | Analyst: <b>DAM</b>   |          |
| Acenaphthene                                     | ND     | 0.092   | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Acenaphthylene                                   | ND     | 0.087   | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Aniline  | ND     | 0.10    | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Anthracene                                       | ND     | 0.071   | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Azobenzene                                       | ND     | 0.13    | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Benz(a)anthracene                                | ND     | 0.092   | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Benzo(a)pyrene                                   | ND     | 0.081   | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Benzo(b)fluoranthene                             | ND     | 0.097   | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Benzo(g,h,i)perylene                             | ND     | 0.095   | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Benzo(k)fluoranthene                             | ND     | 0.095   | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Benzoic acid                                     | ND     | 0.089   | 0.54   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Benzyl alcohol                                   | ND     | 0.084   | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |
| Bis(2-chloroethoxy)methane                       | ND     | 0.12    | 0.21   |      | mg/Kg-dr | 1   | 5/26/2016 6:23:14 PM  | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |



## Hall Environmental Analysis Laboratory, Inc.

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21 (2-2.5')

Project: SWMU 10

Collection Date: 5/18/2016 9:05:00 AM

Lab ID: 1605943-001

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units    | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |      |      |          |    | Analyst: DAM         |          |
| Bis(2-chloroethyl)ether                | ND     | 0.079 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Bis(2-chloroisopropyl)ether            | ND     | 0.096 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Bis(2-ethylhexyl)phthalate             | 0.17   | 0.088 | 0.54 | J    | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 4-Bromophenyl phenyl ether             | ND     | 0.10  | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Butyl benzyl phthalate                 | ND     | 0.095 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Carbazole                              | ND     | 0.073 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 4-Chloro-3-methylphenol                | ND     | 0.13  | 0.54 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 4-Chloroaniline                        | ND     | 0.12  | 0.54 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2-Chloronaphthalene                    | ND     | 0.085 | 0.27 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2-Chlorophenol                         | ND     | 0.085 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 4-Chlorophenyl phenyl ether            | ND     | 0.12  | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Chrysene                               | ND     | 0.091 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Di-n-butyl phthalate                   | ND     | 0.080 | 0.43 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Di-n-octyl phthalate                   | ND     | 0.092 | 0.43 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Dibenz(a,h)anthracene                  | ND     | 0.087 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Dibenzofuran                           | ND     | 0.11  | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 1,2-Dichlorobenzene                    | ND     | 0.082 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 1,3-Dichlorobenzene                    | ND     | 0.083 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 1,4-Dichlorobenzene                    | ND     | 0.091 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 3,3'-Dichlorobenzidine                 | ND     | 0.079 | 0.27 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Diethyl phthalate                      | ND     | 0.11  | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Dimethyl phthalate                     | ND     | 0.11  | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2,4-Dichlorophenol                     | ND     | 0.10  | 0.43 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2,4-Dimethylphenol                     | ND     | 0.12  | 0.32 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 4,6-Dinitro-2-methylphenol             | ND     | 0.065 | 0.43 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2,4-Dinitrophenol                      | ND     | 0.071 | 0.54 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2,4-Dinitrotoluene                     | ND     | 0.096 | 0.54 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2,6-Dinitrotoluene                     | ND     | 0.11  | 0.54 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Fluoranthene                           | ND     | 0.062 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Fluorene                               | ND     | 0.098 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Hexachlorobenzene                      | ND     | 0.085 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Hexachlorobutadiene                    | ND     | 0.12  | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Hexachlorocyclopentadiene              | ND     | 0.12  | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Hexachloroethane                       | ND     | 0.092 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Indeno(1,2,3-cd)pyrene                 | ND     | 0.084 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 1-Methylnaphthalene                    | ND     | 0.11  | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2-Methylnaphthalene                    | ND     | 0.13  | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2-Methylphenol                         | ND     | 0.090 | 0.43 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 3+4-Methylphenol                       | ND     | 0.078 | 0.21 |      | mg/Kg-dr | 1  | 5/26/2016 6:23:14 PM | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-21 (2-2.5')

**Project:** SWMU 10

**Collection Date:** 5/18/2016 9:05:00 AM

**Lab ID:** 1605943-001

**Matrix:** SOIL

**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL      | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|----------|------|-----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |          |      |           |    | Analyst: <b>DAM</b>  |          |
| N-Nitrosodi-n-propylamine              | ND     | 0.10  | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| N-Nitrosodiphenylamine                 | ND     | 0.10  | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Naphthalene                            | ND     | 0.10  | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2-Nitroaniline                         | ND     | 0.12  | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 3-Nitroaniline                         | ND     | 0.095 | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 4-Nitroaniline                         | ND     | 0.076 | 0.43     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Nitrobenzene                           | ND     | 0.11  | 0.43     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2-Nitrophenol                          | ND     | 0.11  | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 4-Nitrophenol                          | ND     | 0.082 | 0.27     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Pentachlorophenol                      | ND     | 0.069 | 0.43     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Phenanthrene                           | ND     | 0.073 | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Phenol                                 | ND     | 0.081 | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Pyrene                                 | ND     | 0.081 | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Pyridine                               | ND     | 0.085 | 0.43     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 1,2,4-Trichlorobenzene                 | ND     | 0.12  | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2,4,5-Trichlorophenol                  | ND     | 0.11  | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| 2,4,6-Trichlorophenol                  | ND     | 0.089 | 0.21     |      | mg/Kg-dr  | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Surr: 2-Fluorophenol                   | 61.9   | 0     | 28.3-102 |      | %Rec      | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Surr: Phenol-d5                        | 65.9   | 0     | 35.7-103 |      | %Rec      | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Surr: 2,4,6-Tribromophenol             | 83.3   | 0     | 35.2-108 |      | %Rec      | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Surr: Nitrobenzene-d5                  | 65.7   |       | 24-118   |      | %Rec      | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Surr: 2-Fluorobiphenyl                 | 80.8   |       | 35.4-111 |      | %Rec      | 1  | 5/26/2016 6:23:14 PM | 25469    |
| Surr: 4-Terphenyl-d14                  | 67.2   |       | 15-91.7  |      | %Rec      | 1  | 5/26/2016 6:23:14 PM | 25469    |
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |          |      |           |    | Analyst: <b>BCN</b>  |          |
| Benzene                                | ND     | 2.24  | 2.24     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Toluene                                | 4.77   | 0.271 | 2.24     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Ethylbenzene                           | 1.22   | 0.301 | 2.24     | J    | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Methyl tert-butyl ether (MTBE)         | ND     | 0.377 | 2.24     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,2,4-Trimethylbenzene                 | 1.95   | 0.380 | 2.24     | J    | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,3,5-Trimethylbenzene                 | 0.774  | 0.370 | 2.24     | J    | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,2-Dichloroethane (EDC)               | ND     | 2.24  | 2.24     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,2-Dibromoethane (EDB)                | ND     | 2.24  | 2.24     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Naphthalene                            | ND     | 2.24  | 2.24     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1-Methylnaphthalene                    | 0.370  | 0.255 | 4.49     | J    | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 2-Methylnaphthalene                    | ND     | 0.590 | 4.49     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Acetone                                | 13.9   | 0.664 | 11.2     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Bromobenzene                           | ND     | 0.228 | 2.24     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Bromodichloromethane                   | ND     | 2.24  | 2.24     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Bromoform                              | ND     | 2.24  | 2.24     |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

**Hall Environmental Analysis Laboratory, Inc.****CLIENT:** Western Refining Company**Client Sample ID:** SWMU 10-21 (2-2.5')**Project:** SWMU 10**Collection Date:** 5/18/2016 9:05:00 AM**Lab ID:** 1605943-001**Matrix:** SOIL**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |      |      |           |    | Analyst: <b>BCN</b>  |          |
| Bromomethane                           | ND     | 0.404 | 3.36 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 2-Butanone                             | ND     | 0.809 | 11.2 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Carbon disulfide                       | ND     | 0.830 | 11.2 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Carbon tetrachloride                   | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Chlorobenzene                          | ND     | 0.257 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Chloroethane                           | ND     | 0.414 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Chloroform                             | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Chloromethane                          | ND     | 0.568 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 2-Chlorotoluene                        | ND     | 0.382 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 4-Chlorotoluene                        | ND     | 0.373 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| cis-1,2-DCE                            | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| cis-1,3-Dichloropropene                | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,2-Dibromo-3-chloropropane            | ND     | 0.239 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Dibromochloromethane                   | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Dibromomethane                         | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,2-Dichlorobenzene                    | ND     | 0.327 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,3-Dichlorobenzene                    | ND     | 0.414 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,4-Dichlorobenzene                    | ND     | 0.420 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Dichlorodifluoromethane                | ND     | 1.31  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,1-Dichloroethane                     | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,1-Dichloroethene                     | ND     | 0.367 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,2-Dichloropropane                    | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,3-Dichloropropane                    | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 2,2-Dichloropropane                    | ND     | 0.283 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,1-Dichloropropene                    | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Hexachlorobutadiene                    | ND     | 0.511 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 2-Hexanone                             | ND     | 0.559 | 11.2 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Isopropylbenzene                       | ND     | 0.282 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 4-Isopropyltoluene                     | ND     | 0.421 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 4-Methyl-2-pentanone                   | ND     | 4.49  | 11.2 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Methylene chloride                     | ND     | 2.24  | 3.36 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| n-Butylbenzene                         | ND     | 0.546 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| n-Propylbenzene                        | ND     | 0.402 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| sec-Butylbenzene                       | ND     | 0.398 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Styrene                                | ND     | 0.287 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| tert-Butylbenzene                      | ND     | 0.328 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,1,1,2-Tetrachloroethane              | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| 1,1,2,2-Tetrachloroethane              | ND     | 2.24  | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |
| Tetrachloroethene (PCE)                | ND     | 0.295 | 2.24 |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21 (2-2.5')

Project: SWMU 10

Collection Date: 5/18/2016 9:05:00 AM

Lab ID: 1605943-001

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                                    | Result | MDL   | PQL    | Qual | Units     | DF | Date Analyzed         | Batch ID |
|---|--------|-------|--------|------|-----------|----|-----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b>      |        |       |        |      |           |    | Analyst: <b>BCN</b>   |          |
| trans-1,2-DCE                               | ND     | 0.227 | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| trans-1,3-Dichloropropene                   | ND     | 0.272 | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| 1,2,3-Trichlorobenzene                      | ND     | 0.543 | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| 1,2,4-Trichlorobenzene                      | ND     | 0.680 | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| 1,1,1-Trichloroethane                       | ND     | 2.24  | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| 1,1,2-Trichloroethane                       | ND     | 2.24  | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| Trichloroethene (TCE)                       | ND     | 2.24  | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| Trichlorofluoromethane                      | ND     | 0.282 | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| 1,2,3-Trichloropropane                      | ND     | 2.24  | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| Vinyl chloride                              | ND     | 0.590 | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| Xylenes, Total                              | 4.46   | 0.894 | 2.24   |      | µg/Kg-dry | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| Surr: 1,2-Dichloroethane-d4                 | 104    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| Surr: 4-Bromofluorobenzene                  | 103    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| Surr: Dibromofluoromethane                  | 93.0   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| Surr: Toluene-d8                            | 97.5   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 4:59:00 PM  | 25468    |
| <b>EPA METHOD 8015D MOD: GASOLINE RANGE</b> |        |       |        |      |           |    | Analyst: <b>DJF</b>   |          |
| Gasoline Range Organics (GRO)               | 2.0    | 0.61  | 4.1    | J    | mg/Kg-dr  | 1  | 5/24/2016 12:36:41 PM | A34454   |
| Surr: BFB                                   | 99.7   | 0     | 70-130 |      | %Rec      | 1  | 5/24/2016 12:36:41 PM | A34454   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21 (12-14')

Project: SWMU 10

Collection Date: 5/18/2016 10:15:00 AM

Lab ID: 1605943-002

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses   | Result | MDL     | PQL    | Qual | Units    | DF  | Date Analyzed         | Batch ID |
|--|--------|---------|--------|------|----------|-----|-----------------------|----------|
| <b>EPA METHOD 8015M/D: DIESEL RANGE ORGANICS</b> |        |         |        |      |          |     | Analyst: <b>KJH</b>   |          |
| Diesel Range Organics (DRO)                      | ND     | 4.5     | 12     |      | mg/Kg-dr | 1   | 5/24/2016 12:43:18 PM | 25449    |
| Motor Oil Range Organics (MRO)                   | ND     | 62      | 62     |      | mg/Kg-dr | 1   | 5/24/2016 12:43:18 PM | 25449    |
| Surr: DNOP                                       | 104    | 0       | 70-130 |      | %Rec     | 1   | 5/24/2016 12:43:18 PM | 25449    |
| <b>PERCENT MOISTURE</b>                          |        |         |        |      |          |     | Analyst: <b>BCN</b>   |          |
| Percent Moisture                                 | 19     | 1.0     | 1.0    |      | wt%      | 1   | 5/23/2016             | R34494   |
| <b>EPA METHOD 7471: MERCURY</b>                  |        |         |        |      |          |     | Analyst: <b>pmf</b>   |          |
| Mercury  | 0.0043 | 0.00069 | 0.040  | J    | mg/Kg-dr | 1   | 5/26/2016 3:46:15 PM  | 25509    |
| <b>EPA METHOD 6010B: SOIL METALS</b>             |        |         |        |      |          |     | Analyst: <b>MED</b>   |          |
| Antimony   | ND     | 7.3     | 15     |      | mg/Kg-dr | 5   | 6/15/2016 6:59:11 AM  | 25508    |
| Arsenic  | 1.8    | 0.90    | 3.1    | J    | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Barium   | 220    | 0.060   | 0.12   |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Beryllium  | 1.3    | 0.0057  | 0.19   |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Cadmium  | ND     | 0.078   | 0.12   |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Chromium   | 11     | 0.15    | 0.37   |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Cobalt   | 5.6    | 0.047   | 0.37   |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Iron   | 20000  | 120     | 310    |      | mg/Kg-dr | 100 | 5/31/2016 10:09:54 AM | 25508    |
| Lead   | 5.1    | 0.22    | 0.31   |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Manganese  | 240    | 0.055   | 0.12   |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Nickel   | 11     | 0.14    | 0.62   |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Selenium   | ND     | 1.3     | 3.1    |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Silver   | ND     | 0.039   | 0.31   |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Vanadium   | 21     | 0.043   | 3.1    |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| Zinc   | 20     | 0.70    | 3.1    |      | mg/Kg-dr | 1   | 5/26/2016 4:08:33 PM  | 25508    |
| <b>EPA METHOD 8270C: SEMIVOLATILES</b>           |        |         |        |      |          |     | Analyst: <b>DAM</b>   |          |
| Acenaphthene                                     | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Acenaphthylene                                   | ND     | 0.10    | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Aniline  | ND     | 0.12    | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Anthracene                                       | ND     | 0.082   | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Azobenzene                                       | ND     | 0.15    | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Benz(a)anthracene                                | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Benzo(a)pyrene                                   | ND     | 0.093   | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Benzo(b)fluoranthene                             | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Benzo(g,h,i)perylene                             | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Benzo(k)fluoranthene                             | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Benzoic acid                                     | ND     | 0.10    | 0.62   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Benzyl alcohol                                   | ND     | 0.097   | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |
| Bis(2-chloroethoxy)methane                       | ND     | 0.13    | 0.25   |      | mg/Kg-dr | 1   | 5/26/2016 6:53:34 PM  | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

## Hall Environmental Analysis Laboratory, Inc.

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21 (12-14')

Project: SWMU 10

Collection Date: 5/18/2016 10:15:00 AM

Lab ID: 1605943-002

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units    | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |      |      |          |    | Analyst: DAM         |          |
| Bis(2-chloroethyl)ether                | ND     | 0.091 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Bis(2-chloroisopropyl)ether            | ND     | 0.11  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Bis(2-ethylhexyl)phthalate             | 0.14   | 0.10  | 0.62 | J    | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 4-Bromophenyl phenyl ether             | ND     | 0.12  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Butyl benzyl phthalate                 | ND     | 0.11  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Carbazole                              | ND     | 0.083 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 4-Chloro-3-methylphenol                | ND     | 0.15  | 0.62 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 4-Chloroaniline                        | ND     | 0.13  | 0.62 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2-Chloronaphthalene                    | ND     | 0.097 | 0.31 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2-Chlorophenol                         | ND     | 0.097 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 4-Chlorophenyl phenyl ether            | ND     | 0.14  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Chrysene                               | ND     | 0.11  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Di-n-butyl phthalate                   | 0.11   | 0.092 | 0.49 | J    | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Di-n-octyl phthalate                   | ND     | 0.11  | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Dibenz(a,h)anthracene                  | ND     | 0.10  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Dibenzofuran                           | ND     | 0.12  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 1,2-Dichlorobenzene                    | ND     | 0.094 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 1,3-Dichlorobenzene                    | ND     | 0.095 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 1,4-Dichlorobenzene                    | ND     | 0.10  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 3,3'-Dichlorobenzidine                 | ND     | 0.091 | 0.31 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Diethyl phthalate                      | ND     | 0.13  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Dimethyl phthalate                     | ND     | 0.12  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2,4-Dichlorophenol                     | ND     | 0.12  | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2,4-Dimethylphenol                     | ND     | 0.13  | 0.37 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 4,6-Dinitro-2-methylphenol             | ND     | 0.075 | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2,4-Dinitrophenol                      | ND     | 0.082 | 0.62 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2,4-Dinitrotoluene                     | ND     | 0.11  | 0.62 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2,6-Dinitrotoluene                     | ND     | 0.13  | 0.62 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Fluoranthene                           | ND     | 0.071 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Fluorene                               | ND     | 0.11  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Hexachlorobenzene                      | ND     | 0.097 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Hexachlorobutadiene                    | ND     | 0.14  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Hexachlorocyclopentadiene              | ND     | 0.14  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Hexachloroethane                       | ND     | 0.11  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Indeno(1,2,3-cd)pyrene                 | ND     | 0.096 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 1-Methylnaphthalene                    | ND     | 0.12  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2-Methylnaphthalene                    | ND     | 0.15  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2-Methylphenol                         | ND     | 0.10  | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 3+4-Methylphenol                       | ND     | 0.089 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 6:53:34 PM | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-21 (12-14')

**Project:** SWMU 10

**Collection Date:** 5/18/2016 10:15:00 AM

**Lab ID:** 1605943-002

**Matrix:** SOIL

**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL      | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|----------|------|-----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |          |      |           |    | Analyst: <b>DAM</b>  |          |
| N-Nitrosodi-n-propylamine              | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| N-Nitrosodiphenylamine                 | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Naphthalene                            | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2-Nitroaniline                         | ND     | 0.13  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 3-Nitroaniline                         | ND     | 0.11  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 4-Nitroaniline                         | ND     | 0.087 | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Nitrobenzene                           | ND     | 0.13  | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2-Nitrophenol                          | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 4-Nitrophenol                          | ND     | 0.094 | 0.31     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Pentachlorophenol                      | ND     | 0.079 | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Phenanthrene                           | ND     | 0.084 | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Phenol                                 | ND     | 0.093 | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Pyrene                                 | ND     | 0.093 | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Pyridine                               | ND     | 0.098 | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 1,2,4-Trichlorobenzene                 | ND     | 0.13  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2,4,5-Trichlorophenol                  | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| 2,4,6-Trichlorophenol                  | ND     | 0.10  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Surr: 2-Fluorophenol                   | 48.1   | 0     | 28.3-102 |      | %Rec      | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Surr: Phenol-d5                        | 55.6   | 0     | 35.7-103 |      | %Rec      | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Surr: 2,4,6-Tribromophenol             | 57.9   | 0     | 35.2-108 |      | %Rec      | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Surr: Nitrobenzene-d5                  | 57.3   |       | 24-118   |      | %Rec      | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Surr: 2-Fluorobiphenyl                 | 57.6   |       | 35.4-111 |      | %Rec      | 1  | 5/26/2016 6:53:34 PM | 25469    |
| Surr: 4-Terphenyl-d14                  | 56.0   |       | 15-91.7  |      | %Rec      | 1  | 5/26/2016 6:53:34 PM | 25469    |
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |          |      |           |    | Analyst: <b>BCN</b>  |          |
| Benzene                                | 1.93   | 1.87  | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Toluene                                | 1.47   | 0.226 | 1.87     | J    | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Ethylbenzene                           | ND     | 0.252 | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Methyl tert-butyl ether (MTBE)         | ND     | 0.315 | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,2,4-Trimethylbenzene                 | 0.347  | 0.318 | 1.87     | J    | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,3,5-Trimethylbenzene                 | ND     | 0.309 | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,2-Dichloroethane (EDC)               | ND     | 1.87  | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,2-Dibromoethane (EDB)                | ND     | 1.87  | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Naphthalene                            | ND     | 1.87  | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1-Methylnaphthalene                    | ND     | 0.213 | 3.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 2-Methylnaphthalene                    | ND     | 0.493 | 3.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Acetone                                | 5.84   | 0.554 | 9.37     | J    | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Bromobenzene                           | ND     | 0.191 | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Bromodichloromethane                   | ND     | 1.87  | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Bromoform                              | ND     | 1.87  | 1.87     |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

**Hall Environmental Analysis Laboratory, Inc.****CLIENT:** Western Refining Company**Client Sample ID:** SWMU 10-21 (12-14')**Project:** SWMU 10**Collection Date:** 5/18/2016 10:15:00 AM**Lab ID:** 1605943-002**Matrix:** SOIL**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |      |      |           |    | Analyst: <b>BCN</b>  |          |
| Bromomethane                           | ND     | 0.337 | 2.81 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 2-Butanone                             | ND     | 0.675 | 9.37 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Carbon disulfide                       | ND     | 0.693 | 9.37 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Carbon tetrachloride                   | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Chlorobenzene                          | ND     | 0.215 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Chloroethane                           | ND     | 0.346 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Chloroform                             | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Chloromethane                          | ND     | 0.475 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 2-Chlorotoluene                        | ND     | 0.319 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 4-Chlorotoluene                        | ND     | 0.311 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| cis-1,2-DCE                            | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| cis-1,3-Dichloropropene                | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,2-Dibromo-3-chloropropane            | ND     | 0.199 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Dibromochloromethane                   | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Dibromomethane                         | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,2-Dichlorobenzene                    | ND     | 0.273 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,3-Dichlorobenzene                    | ND     | 0.346 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,4-Dichlorobenzene                    | ND     | 0.351 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Dichlorodifluoromethane                | ND     | 1.09  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,1-Dichloroethane                     | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,1-Dichloroethene                     | ND     | 0.307 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,2-Dichloropropane                    | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,3-Dichloropropane                    | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 2,2-Dichloropropane                    | ND     | 0.237 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,1-Dichloropropene                    | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Hexachlorobutadiene                    | ND     | 0.427 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 2-Hexanone                             | ND     | 0.467 | 9.37 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Isopropylbenzene                       | ND     | 0.236 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 4-Isopropyltoluene                     | ND     | 0.351 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 4-Methyl-2-pentanone                   | ND     | 3.75  | 9.37 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Methylene chloride                     | ND     | 1.87  | 2.81 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| n-Butylbenzene                         | ND     | 0.456 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| n-Propylbenzene                        | ND     | 0.335 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| sec-Butylbenzene                       | ND     | 0.333 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Styrene                                | ND     | 0.239 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| tert-Butylbenzene                      | ND     | 0.274 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,1,1,2-Tetrachloroethane              | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,1,2,2-Tetrachloroethane              | ND     | 1.87  | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Tetrachloroethene (PCE)                | ND     | 0.246 | 1.87 |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |



# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21 (12-14')

Project: SWMU 10

Collection Date: 5/18/2016 10:15:00 AM

Lab ID: 1605943-002

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                                    | Result | MDL   | PQL    | Qual | Units     | DF | Date Analyzed        | Batch ID |
|---|--------|-------|--------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b>      |        |       |        |      |           |    | Analyst: <b>BCN</b>  |          |
| trans-1,2-DCE                               | ND     | 0.189 | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| trans-1,3-Dichloropropene                   | ND     | 0.227 | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,2,3-Trichlorobenzene                      | ND     | 0.453 | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,2,4-Trichlorobenzene                      | ND     | 0.568 | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,1,1-Trichloroethane                       | ND     | 1.87  | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,1,2-Trichloroethane                       | ND     | 1.87  | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Trichloroethene (TCE)                       | ND     | 1.87  | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Trichlorofluoromethane                      | ND     | 0.235 | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| 1,2,3-Trichloropropane                      | ND     | 1.87  | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Vinyl chloride                              | ND     | 0.493 | 1.87   |      | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Xylenes, Total                              | 0.937  | 0.747 | 1.87   | J    | µg/Kg-dry | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Surr: 1,2-Dichloroethane-d4                 | 99.7   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Surr: 4-Bromofluorobenzene                  | 102    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Surr: Dibromofluoromethane                  | 91.9   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 5:24:00 PM | 25468    |
| Surr: Toluene-d8                            | 97.2   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 5:24:00 PM | 25468    |
| <b>EPA METHOD 8015D MOD: GASOLINE RANGE</b> |        |       |        |      |           |    | Analyst: <b>DJF</b>  |          |
| Gasoline Range Organics (GRO)               | 0.60   | 0.48  | 3.2    | J    | mg/Kg-dr  | 1  | 5/24/2016 1:05:04 PM | A34454   |
| Surr: BFB                                   | 93.4   | 0     | 70-130 |      | %Rec      | 1  | 5/24/2016 1:05:04 PM | A34454   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21 (20-22')

Project: SWMU 10

Collection Date: 5/18/2016 10:25:00 AM

Lab ID: 1605943-003

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses   | Result | MDL     | PQL    | Qual | Units    | DF  | Date Analyzed         | Batch ID |
|--|--------|---------|--------|------|----------|-----|-----------------------|----------|
| <b>EPA METHOD 8015M/D: DIESEL RANGE ORGANICS</b> |        |         |        |      |          |     | Analyst: <b>KJH</b>   |          |
| Diesel Range Organics (DRO)                      | ND     | 3.8     | 11     |      | mg/Kg-dr | 1   | 5/24/2016 1:06:24 PM  | 25449    |
| Motor Oil Range Organics (MRO)                   | ND     | 53      | 53     |      | mg/Kg-dr | 1   | 5/24/2016 1:06:24 PM  | 25449    |
| Surr: DNOP                                       | 120    | 0       | 70-130 |      | %Rec     | 1   | 5/24/2016 1:06:24 PM  | 25449    |
| <b>PERCENT MOISTURE</b>                          |        |         |        |      |          |     | Analyst: <b>BCN</b>   |          |
| Percent Moisture                                 | 12     | 1.0     | 1.0    |      | wt%      | 1   | 5/23/2016             | R34494   |
| <b>EPA METHOD 7471: MERCURY</b>                  |        |         |        |      |          |     | Analyst: <b>pmf</b>   |          |
| Mercury  | ND     | 0.00063 | 0.037  |      | mg/Kg-dr | 1   | 5/26/2016 3:48:01 PM  | 25509    |
| <b>EPA METHOD 6010B: SOIL METALS</b>             |        |         |        |      |          |     | Analyst: <b>MED</b>   |          |
| Antimony   | ND     | 6.8     | 14     |      | mg/Kg-dr | 5   | 5/31/2016 10:12:55 AM | 25508    |
| Arsenic  | 3.3    | 0.83    | 2.9    |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Barium   | 1000   | 0.28    | 0.57   |      | mg/Kg-dr | 5   | 5/31/2016 10:12:55 AM | 25508    |
| Beryllium  | 0.91   | 0.0053  | 0.17   |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Cadmium  | ND     | 0.072   | 0.11   |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Chromium   | 12     | 0.14    | 0.34   |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Cobalt   | 8.8    | 0.043   | 0.34   |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Iron   | 18000  | 110     | 290    |      | mg/Kg-dr | 100 | 5/31/2016 10:11:24 AM | 25508    |
| Lead   | 2.6    | 0.20    | 0.29   |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Manganese  | 1800   | 0.51    | 1.1    |      | mg/Kg-dr | 10  | 5/31/2016 11:01:26 AM | 25508    |
| Nickel   | 15     | 0.13    | 0.57   |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Selenium   | ND     | 1.2     | 2.9    |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Silver   | ND     | 0.036   | 0.29   |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Vanadium   | 17     | 0.040   | 2.9    |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| Zinc   | 27     | 0.65    | 2.9    |      | mg/Kg-dr | 1   | 5/26/2016 4:10:03 PM  | 25508    |
| <b>EPA METHOD 8270C: SEMIVOLATILES</b>           |        |         |        |      |          |     | Analyst: <b>DAM</b>   |          |
| Acenaphthene                                     | ND     | 0.096   | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Acenaphthylene                                   | ND     | 0.091   | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Aniline  | ND     | 0.11    | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Anthracene                                       | ND     | 0.074   | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Azobenzene                                       | ND     | 0.14    | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Benz(a)anthracene                                | ND     | 0.096   | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Benzo(a)pyrene                                   | ND     | 0.085   | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Benzo(b)fluoranthene                             | ND     | 0.10    | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Benzo(g,h,i)perylene                             | ND     | 0.099   | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Benzo(k)fluoranthene                             | ND     | 0.099   | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Benzoic acid                                     | ND     | 0.093   | 0.56   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Benzyl alcohol                                   | ND     | 0.088   | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |
| Bis(2-chloroethoxy)methane                       | ND     | 0.12    | 0.22   |      | mg/Kg-dr | 1   | 5/26/2016 7:23:47 PM  | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

## Hall Environmental Analysis Laboratory, Inc.

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21 (20-22')

Project: SWMU 10

Collection Date: 5/18/2016 10:25:00 AM

Lab ID: 1605943-003

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units    | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |      |      |          |    | Analyst: DAM         |          |
| Bis(2-chloroethyl)ether                | ND     | 0.082 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Bis(2-chloroisopropyl)ether            | ND     | 0.10  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Bis(2-ethylhexyl)phthalate             | 0.12   | 0.091 | 0.56 | J    | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 4-Bromophenyl phenyl ether             | ND     | 0.11  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Butyl benzyl phthalate                 | ND     | 0.099 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Carbazole                              | ND     | 0.076 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 4-Chloro-3-methylphenol                | ND     | 0.13  | 0.56 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 4-Chloroaniline                        | ND     | 0.12  | 0.56 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2-Chloronaphthalene                    | ND     | 0.088 | 0.28 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2-Chlorophenol                         | ND     | 0.088 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 4-Chlorophenyl phenyl ether            | ND     | 0.13  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Chrysene                               | ND     | 0.095 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Di-n-butyl phthalate                   | ND     | 0.084 | 0.45 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Di-n-octyl phthalate                   | ND     | 0.096 | 0.45 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Dibenz(a,h)anthracene                  | ND     | 0.091 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Dibenzofuran                           | ND     | 0.11  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 1,2-Dichlorobenzene                    | ND     | 0.086 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 1,3-Dichlorobenzene                    | ND     | 0.087 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 1,4-Dichlorobenzene                    | ND     | 0.095 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 3,3'-Dichlorobenzidine                 | ND     | 0.083 | 0.28 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Diethyl phthalate                      | ND     | 0.11  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Dimethyl phthalate                     | ND     | 0.11  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2,4-Dichlorophenol                     | ND     | 0.10  | 0.45 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2,4-Dimethylphenol                     | ND     | 0.12  | 0.34 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 4,6-Dinitro-2-methylphenol             | ND     | 0.068 | 0.45 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2,4-Dinitrophenol                      | ND     | 0.074 | 0.56 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2,4-Dinitrotoluene                     | ND     | 0.10  | 0.56 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2,6-Dinitrotoluene                     | ND     | 0.12  | 0.56 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Fluoranthene                           | ND     | 0.065 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Fluorene                               | ND     | 0.10  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Hexachlorobenzene                      | ND     | 0.088 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Hexachlorobutadiene                    | ND     | 0.13  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Hexachlorocyclopentadiene              | ND     | 0.13  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Hexachloroethane                       | ND     | 0.096 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Indeno(1,2,3-cd)pyrene                 | ND     | 0.087 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 1-Methylnaphthalene                    | ND     | 0.11  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2-Methylnaphthalene                    | ND     | 0.13  | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2-Methylphenol                         | ND     | 0.094 | 0.45 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 3+4-Methylphenol                       | ND     | 0.081 | 0.22 |      | mg/Kg-dr | 1  | 5/26/2016 7:23:47 PM | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-21 (20-22')

**Project:** SWMU 10

**Collection Date:** 5/18/2016 10:25:00 AM

**Lab ID:** 1605943-003

**Matrix:** SOIL

**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL      | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|----------|------|-----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |          |      |           |    | Analyst: <b>DAM</b>  |          |
| N-Nitrosodi-n-propylamine              | ND     | 0.11  | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| N-Nitrosodiphenylamine                 | ND     | 0.11  | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Naphthalene                            | ND     | 0.11  | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2-Nitroaniline                         | ND     | 0.12  | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 3-Nitroaniline                         | ND     | 0.099 | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 4-Nitroaniline                         | ND     | 0.079 | 0.45     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Nitrobenzene                           | ND     | 0.12  | 0.45     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2-Nitrophenol                          | ND     | 0.11  | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 4-Nitrophenol                          | ND     | 0.085 | 0.28     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Pentachlorophenol                      | ND     | 0.072 | 0.45     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Phenanthrene                           | ND     | 0.076 | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Phenol                                 | ND     | 0.084 | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Pyrene                                 | ND     | 0.085 | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Pyridine                               | ND     | 0.089 | 0.45     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 1,2,4-Trichlorobenzene                 | ND     | 0.12  | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2,4,5-Trichlorophenol                  | ND     | 0.11  | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| 2,4,6-Trichlorophenol                  | ND     | 0.093 | 0.22     |      | mg/Kg-dr  | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Surr: 2-Fluorophenol                   | 50.2   | 0     | 28.3-102 |      | %Rec      | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Surr: Phenol-d5                        | 54.9   | 0     | 35.7-103 |      | %Rec      | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Surr: 2,4,6-Tribromophenol             | 67.1   | 0     | 35.2-108 |      | %Rec      | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Surr: Nitrobenzene-d5                  | 52.9   |       | 24-118   |      | %Rec      | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Surr: 2-Fluorobiphenyl                 | 64.6   |       | 35.4-111 |      | %Rec      | 1  | 5/26/2016 7:23:47 PM | 25469    |
| Surr: 4-Terphenyl-d14                  | 51.3   |       | 15-91.7  |      | %Rec      | 1  | 5/26/2016 7:23:47 PM | 25469    |
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |          |      |           |    | Analyst: <b>BCN</b>  |          |
| Benzene                                | 8.48   | 1.75  | 1.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Toluene                                | 6.22   | 0.212 | 1.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Ethylbenzene                           | 1.32   | 0.236 | 1.75     | J    | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Methyl tert-butyl ether (MTBE)         | ND     | 0.295 | 1.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,2,4-Trimethylbenzene                 | 1.03   | 0.297 | 1.75     | J    | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,3,5-Trimethylbenzene                 | 0.684  | 0.289 | 1.75     | J    | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,2-Dichloroethane (EDC)               | ND     | 1.75  | 1.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,2-Dibromoethane (EDB)                | ND     | 1.75  | 1.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Naphthalene                            | ND     | 1.75  | 1.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1-Methylnaphthalene                    | 0.219  | 0.199 | 3.51     | J    | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 2-Methylnaphthalene                    | ND     | 0.462 | 3.51     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Acetone                                | 6.34   | 0.519 | 8.77     | J    | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Bromobenzene                           | ND     | 0.178 | 1.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Bromodichloromethane                   | ND     | 1.75  | 1.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Bromoform                              | ND     | 1.75  | 1.75     |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

**Hall Environmental Analysis Laboratory, Inc.****CLIENT:** Western Refining Company**Client Sample ID:** SWMU 10-21 (20-22')**Project:** SWMU 10**Collection Date:** 5/18/2016 10:25:00 AM**Lab ID:** 1605943-003**Matrix:** SOIL**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |      |      |           |    | Analyst: <b>BCN</b>  |          |
| Bromomethane                           | ND     | 0.316 | 2.63 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 2-Butanone                             | ND     | 0.632 | 8.77 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Carbon disulfide                       | ND     | 0.649 | 8.77 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Carbon tetrachloride                   | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Chlorobenzene                          | ND     | 0.201 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Chloroethane                           | ND     | 0.324 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Chloroform                             | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Chloromethane                          | ND     | 0.444 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 2-Chlorotoluene                        | ND     | 0.298 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 4-Chlorotoluene                        | ND     | 0.292 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| cis-1,2-DCE                            | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| cis-1,3-Dichloropropene                | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,2-Dibromo-3-chloropropane            | ND     | 0.187 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Dibromochloromethane                   | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Dibromomethane                         | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,2-Dichlorobenzene                    | ND     | 0.255 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,3-Dichlorobenzene                    | ND     | 0.324 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,4-Dichlorobenzene                    | ND     | 0.328 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Dichlorodifluoromethane                | ND     | 1.02  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,1-Dichloroethane                     | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,1-Dichloroethene                     | ND     | 0.287 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,2-Dichloropropane                    | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,3-Dichloropropane                    | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 2,2-Dichloropropane                    | ND     | 0.221 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,1-Dichloropropene                    | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Hexachlorobutadiene                    | ND     | 0.400 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 2-Hexanone                             | ND     | 0.437 | 8.77 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Isopropylbenzene                       | ND     | 0.221 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 4-Isopropyltoluene                     | ND     | 0.329 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 4-Methyl-2-pentanone                   | ND     | 3.51  | 8.77 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Methylene chloride                     | ND     | 1.75  | 2.63 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| n-Butylbenzene                         | ND     | 0.427 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| n-Propylbenzene                        | ND     | 0.314 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| sec-Butylbenzene                       | ND     | 0.311 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Styrene                                | ND     | 0.224 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| tert-Butylbenzene                      | ND     | 0.256 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,1,1,2-Tetrachloroethane              | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,1,2,2-Tetrachloroethane              | ND     | 1.75  | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Tetrachloroethene (PCE)                | ND     | 0.230 | 1.75 |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21 (20-22')

Project: SWMU 10

Collection Date: 5/18/2016 10:25:00 AM

Lab ID: 1605943-003

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                                    | Result | MDL   | PQL    | Qual | Units     | DF | Date Analyzed        | Batch ID |
|---|--------|-------|--------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b>      |        |       |        |      |           |    | Analyst: <b>BCN</b>  |          |
| trans-1,2-DCE                               | ND     | 0.177 | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| trans-1,3-Dichloropropene                   | ND     | 0.212 | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,2,3-Trichlorobenzene                      | ND     | 0.424 | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,2,4-Trichlorobenzene                      | ND     | 0.532 | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,1,1-Trichloroethane                       | ND     | 1.75  | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,1,2-Trichloroethane                       | ND     | 1.75  | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Trichloroethene (TCE)                       | ND     | 1.75  | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Trichlorofluoromethane                      | ND     | 0.220 | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| 1,2,3-Trichloropropane                      | ND     | 1.75  | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Vinyl chloride                              | ND     | 0.461 | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Xylenes, Total                              | 3.85   | 0.699 | 1.75   |      | µg/Kg-dry | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Surr: 1,2-Dichloroethane-d4                 | 106    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Surr: 4-Bromofluorobenzene                  | 101    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Surr: Dibromofluoromethane                  | 92.3   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 5:49:00 PM | 25468    |
| Surr: Toluene-d8                            | 96.1   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 5:49:00 PM | 25468    |
| <b>EPA METHOD 8015D MOD: GASOLINE RANGE</b> |        |       |        |      |           |    | Analyst: <b>DJF</b>  |          |
| Gasoline Range Organics (GRO)               | 0.76   | 0.54  | 3.6    | J    | mg/Kg-dr  | 1  | 5/24/2016 1:33:30 PM | A34454   |
| Surr: BFB                                   | 101    | 0     | 70-130 |      | %Rec      | 1  | 5/24/2016 1:33:30 PM | A34454   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-22 (2-2.5')

Project: SWMU 10

Collection Date: 5/18/2016 4:35:00 PM

Lab ID: 1605943-004

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses   | Result | MDL     | PQL    | Qual | Units    | DF | Date Analyzed         | Batch ID |
|--|--------|---------|--------|------|----------|----|-----------------------|----------|
| <b>EPA METHOD 8015M/D: DIESEL RANGE ORGANICS</b> |        |         |        |      |          |    | Analyst: <b>KJH</b>   |          |
| Diesel Range Organics (DRO)                      | ND     | 4.5     | 12     |      | mg/Kg-dr | 1  | 5/24/2016 1:28:09 PM  | 25449    |
| Motor Oil Range Organics (MRO)                   | ND     | 62      | 62     |      | mg/Kg-dr | 1  | 5/24/2016 1:28:09 PM  | 25449    |
| Surr: DNOP                                       | 109    | 0       | 70-130 |      | %Rec     | 1  | 5/24/2016 1:28:09 PM  | 25449    |
| <b>PERCENT MOISTURE</b>                          |        |         |        |      |          |    | Analyst: <b>BCN</b>   |          |
| Percent Moisture                                 | 19     | 1.0     | 1.0    |      | wt%      | 1  | 5/23/2016             | R34494   |
| <b>EPA METHOD 7471: MERCURY</b>                  |        |         |        |      |          |    | Analyst: <b>pmf</b>   |          |
| Mercury  | 0.044  | 0.00071 | 0.041  |      | mg/Kg-dr | 1  | 5/26/2016 3:49:49 PM  | 25509    |
| <b>EPA METHOD 6010B: SOIL METALS</b>             |        |         |        |      |          |    | Analyst: <b>MED</b>   |          |
| Antimony   | ND     | 7.3     | 15     |      | mg/Kg-dr | 5  | 5/31/2016 10:16:13 AM | 25508    |
| Arsenic  | 2.4    | 0.90    | 3.1    | J    | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Barium   | 290    | 0.060   | 0.12   |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Beryllium  | 0.89   | 0.0057  | 0.18   |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Cadmium  | ND     | 0.078   | 0.12   |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Chromium   | 12     | 0.15    | 0.37   |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Cobalt   | 4.9    | 0.047   | 0.37   |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Iron   | 14000  | 60      | 150    |      | mg/Kg-dr | 50 | 5/31/2016 10:14:41 AM | 25508    |
| Lead   | 4.3    | 0.22    | 0.31   |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Manganese  | 570    | 0.27    | 0.62   |      | mg/Kg-dr | 5  | 5/31/2016 10:16:13 AM | 25508    |
| Nickel   | 8.6    | 0.14    | 0.62   |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Selenium   | ND     | 1.3     | 3.1    |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Silver   | ND     | 0.039   | 0.31   |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Vanadium   | 18     | 0.043   | 3.1    |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| Zinc   | 20     | 0.70    | 3.1    |      | mg/Kg-dr | 1  | 5/26/2016 4:11:42 PM  | 25508    |
| <b>EPA METHOD 8270C: SEMIVOLATILES</b>           |        |         |        |      |          |    | Analyst: <b>DAM</b>   |          |
| Acenaphthene                                     | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Acenaphthylene                                   | ND     | 0.10    | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Aniline  | ND     | 0.12    | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Anthracene                                       | ND     | 0.081   | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Azobenzene                                       | ND     | 0.15    | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Benz(a)anthracene                                | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Benzo(a)pyrene                                   | ND     | 0.093   | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Benzo(b)fluoranthene                             | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Benzo(g,h,i)perylene                             | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Benzo(k)fluoranthene                             | ND     | 0.11    | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Benzoic acid                                     | ND     | 0.10    | 0.61   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Benzyl alcohol                                   | ND     | 0.096   | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |
| Bis(2-chloroethoxy)methane                       | ND     | 0.13    | 0.25   |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM  | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

## Hall Environmental Analysis Laboratory, Inc.

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-22 (2-2.5')

Project: SWMU 10

Collection Date: 5/18/2016 4:35:00 PM

Lab ID: 1605943-004

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units    | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |      |      |          |    | Analyst: DAM         |          |
| Bis(2-chloroethyl)ether                | ND     | 0.090 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Bis(2-chloroisopropyl)ether            | ND     | 0.11  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Bis(2-ethylhexyl)phthalate             | 0.19   | 0.10  | 0.61 | J    | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 4-Bromophenyl phenyl ether             | ND     | 0.12  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Butyl benzyl phthalate                 | ND     | 0.11  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Carbazole                              | ND     | 0.083 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 4-Chloro-3-methylphenol                | ND     | 0.15  | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 4-Chloroaniline                        | ND     | 0.13  | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2-Chloronaphthalene                    | ND     | 0.097 | 0.31 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2-Chlorophenol                         | ND     | 0.097 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 4-Chlorophenyl phenyl ether            | ND     | 0.14  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Chrysene                               | ND     | 0.10  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Di-n-butyl phthalate                   | 0.12   | 0.092 | 0.49 | J    | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Di-n-octyl phthalate                   | ND     | 0.10  | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Dibenz(a,h)anthracene                  | ND     | 0.099 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Dibenzofuran                           | ND     | 0.12  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 1,2-Dichlorobenzene                    | ND     | 0.094 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 1,3-Dichlorobenzene                    | ND     | 0.095 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 1,4-Dichlorobenzene                    | ND     | 0.10  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 3,3'-Dichlorobenzidine                 | ND     | 0.090 | 0.31 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Diethyl phthalate                      | ND     | 0.12  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Dimethyl phthalate                     | ND     | 0.12  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2,4-Dichlorophenol                     | ND     | 0.11  | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2,4-Dimethylphenol                     | ND     | 0.13  | 0.37 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 4,6-Dinitro-2-methylphenol             | ND     | 0.074 | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2,4-Dinitrophenol                      | ND     | 0.081 | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2,4-Dinitrotoluene                     | ND     | 0.11  | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2,6-Dinitrotoluene                     | ND     | 0.13  | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Fluoranthene                           | ND     | 0.071 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Fluorene                               | ND     | 0.11  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Hexachlorobenzene                      | ND     | 0.097 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Hexachlorobutadiene                    | ND     | 0.14  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Hexachlorocyclopentadiene              | ND     | 0.14  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Hexachloroethane                       | ND     | 0.11  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Indeno(1,2,3-cd)pyrene                 | ND     | 0.096 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 1-Methylnaphthalene                    | ND     | 0.12  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2-Methylnaphthalene                    | ND     | 0.15  | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2-Methylphenol                         | ND     | 0.10  | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 3+4-Methylphenol                       | ND     | 0.089 | 0.25 |      | mg/Kg-dr | 1  | 5/26/2016 7:53:55 PM | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |



# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-22 (2-2.5')

**Project:** SWMU 10

**Collection Date:** 5/18/2016 4:35:00 PM

**Lab ID:** 1605943-004

**Matrix:** SOIL

**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL      | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|----------|------|-----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |          |      |           |    | Analyst: <b>DAM</b>  |          |
| N-Nitrosodi-n-propylamine              | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| N-Nitrosodiphenylamine                 | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Naphthalene                            | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2-Nitroaniline                         | ND     | 0.13  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 3-Nitroaniline                         | ND     | 0.11  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 4-Nitroaniline                         | ND     | 0.086 | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Nitrobenzene                           | ND     | 0.13  | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2-Nitrophenol                          | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 4-Nitrophenol                          | ND     | 0.093 | 0.31     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Pentachlorophenol                      | ND     | 0.079 | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Phenanthrene                           | ND     | 0.083 | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Phenol                                 | ND     | 0.092 | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Pyrene                                 | ND     | 0.093 | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Pyridine                               | ND     | 0.097 | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 1,2,4-Trichlorobenzene                 | ND     | 0.13  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2,4,5-Trichlorophenol                  | ND     | 0.12  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| 2,4,6-Trichlorophenol                  | ND     | 0.10  | 0.25     |      | mg/Kg-dr  | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Surr: 2-Fluorophenol                   | 62.9   | 0     | 28.3-102 |      | %Rec      | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Surr: Phenol-d5                        | 64.7   | 0     | 35.7-103 |      | %Rec      | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Surr: 2,4,6-Tribromophenol             | 83.2   | 0     | 35.2-108 |      | %Rec      | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Surr: Nitrobenzene-d5                  | 66.7   |       | 24-118   |      | %Rec      | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Surr: 2-Fluorobiphenyl                 | 70.6   |       | 35.4-111 |      | %Rec      | 1  | 5/26/2016 7:53:55 PM | 25469    |
| Surr: 4-Terphenyl-d14                  | 95.9   |       | 15-91.7  | S    | %Rec      | 1  | 5/26/2016 7:53:55 PM | 25469    |
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |          |      |           |    | Analyst: <b>BCN</b>  |          |
| Benzene                                | ND     | 1.67  | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Toluene                                | 1.18   | 0.201 | 1.67     | J    | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Ethylbenzene                           | ND     | 0.224 | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Methyl tert-butyl ether (MTBE)         | ND     | 0.280 | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,2,4-Trimethylbenzene                 | 0.300  | 0.283 | 1.67     | J    | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,3,5-Trimethylbenzene                 | ND     | 0.275 | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,2-Dichloroethane (EDC)               | ND     | 1.67  | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,2-Dibromoethane (EDB)                | ND     | 1.67  | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Naphthalene                            | ND     | 1.67  | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1-Methylnaphthalene                    | ND     | 0.189 | 3.33     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 2-Methylnaphthalene                    | ND     | 0.439 | 3.33     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Acetone                                | 11.7   | 0.493 | 8.34     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Bromobenzene                           | ND     | 0.170 | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Bromodichloromethane                   | ND     | 1.67  | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Bromoform                              | ND     | 1.67  | 1.67     |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-22 (2-2.5')

**Project:** SWMU 10

**Collection Date:** 5/18/2016 4:35:00 PM

**Lab ID:** 1605943-004

**Matrix:** SOIL

**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |      |      |           |    | Analyst: <b>BCN</b>  |          |
| Bromomethane                           | ND     | 0.300 | 2.50 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 2-Butanone                             | ND     | 0.601 | 8.34 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Carbon disulfide                       | ND     | 0.617 | 8.34 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Carbon tetrachloride                   | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Chlorobenzene                          | ND     | 0.191 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Chloroethane                           | ND     | 0.308 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Chloroform                             | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Chloromethane                          | ND     | 0.422 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 2-Chlorotoluene                        | ND     | 0.284 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 4-Chlorotoluene                        | ND     | 0.277 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| cis-1,2-DCE                            | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| cis-1,3-Dichloropropene                | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,2-Dibromo-3-chloropropane            | ND     | 0.177 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Dibromochloromethane                   | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Dibromomethane                         | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,2-Dichlorobenzene                    | ND     | 0.243 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,3-Dichlorobenzene                    | ND     | 0.308 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,4-Dichlorobenzene                    | ND     | 0.312 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Dichlorodifluoromethane                | ND     | 0.973 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,1-Dichloroethane                     | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,1-Dichloroethene                     | ND     | 0.273 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,2-Dichloropropane                    | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,3-Dichloropropane                    | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 2,2-Dichloropropane                    | ND     | 0.210 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,1-Dichloropropene                    | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Hexachlorobutadiene                    | ND     | 0.380 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 2-Hexanone                             | ND     | 0.415 | 8.34 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Isopropylbenzene                       | ND     | 0.210 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 4-Isopropyltoluene                     | ND     | 0.313 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 4-Methyl-2-pentanone                   | ND     | 3.33  | 8.34 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Methylene chloride                     | ND     | 1.67  | 2.50 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| n-Butylbenzene                         | ND     | 0.406 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| n-Propylbenzene                        | ND     | 0.298 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| sec-Butylbenzene                       | ND     | 0.296 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Styrene                                | ND     | 0.213 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| tert-Butylbenzene                      | ND     | 0.244 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,1,1,2-Tetrachloroethane              | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,1,2,2-Tetrachloroethane              | ND     | 1.67  | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Tetrachloroethene (PCE)                | ND     | 0.219 | 1.67 |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-22 (2-2.5')

Project: SWMU 10

Collection Date: 5/18/2016 4:35:00 PM

Lab ID: 1605943-004

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                                    | Result | MDL   | PQL    | Qual | Units     | DF | Date Analyzed        | Batch ID |
|---|--------|-------|--------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b>      |        |       |        |      |           |    | Analyst: <b>BCN</b>  |          |
| trans-1,2-DCE                               | ND     | 0.168 | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| trans-1,3-Dichloropropene                   | ND     | 0.202 | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,2,3-Trichlorobenzene                      | ND     | 0.403 | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,2,4-Trichlorobenzene                      | ND     | 0.506 | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,1,1-Trichloroethane                       | ND     | 1.67  | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,1,2-Trichloroethane                       | ND     | 1.67  | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Trichloroethene (TCE)                       | ND     | 1.67  | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Trichlorofluoromethane                      | ND     | 0.209 | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| 1,2,3-Trichloropropane                      | ND     | 1.67  | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Vinyl chloride                              | ND     | 0.439 | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Xylenes, Total                              | ND     | 0.664 | 1.67   |      | µg/Kg-dry | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Surr: 1,2-Dichloroethane-d4                 | 104    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Surr: 4-Bromofluorobenzene                  | 102    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Surr: Dibromofluoromethane                  | 93.8   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 6:14:00 PM | 25468    |
| Surr: Toluene-d8                            | 97.4   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 6:14:00 PM | 25468    |
| <b>EPA METHOD 8015D MOD: GASOLINE RANGE</b> |        |       |        |      |           |    | Analyst: <b>DJF</b>  |          |
| Gasoline Range Organics (GRO)               | 0.60   | 0.49  | 3.2    | J    | mg/Kg-dr  | 1  | 5/24/2016 2:01:48 PM | A34454   |
| Surr: BFB                                   | 97.6   | 0     | 70-130 |      | %Rec      | 1  | 5/24/2016 2:01:48 PM | A34454   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-22 (8-9')

Project: SWMU 10

Collection Date: 5/18/2016 4:50:00 PM

Lab ID: 1605943-005

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses   | Result | MDL     | PQL    | Qual | Units    | DF  | Date Analyzed         | Batch ID |
|--|--------|---------|--------|------|----------|-----|-----------------------|----------|
| <b>EPA METHOD 8015M/D: DIESEL RANGE ORGANICS</b> |        |         |        |      |          |     | Analyst: <b>KJH</b>   |          |
| Diesel Range Organics (DRO)                      | ND     | 4.1     | 11     |      | mg/Kg-dr | 1   | 5/24/2016 2:11:34 PM  | 25449    |
| Motor Oil Range Organics (MRO)                   | ND     | 57      | 57     |      | mg/Kg-dr | 1   | 5/24/2016 2:11:34 PM  | 25449    |
| Surr: DNOP                                       | 92.7   | 0       | 70-130 |      | %Rec     | 1   | 5/24/2016 2:11:34 PM  | 25449    |
| <b>PERCENT MOISTURE</b>                          |        |         |        |      |          |     | Analyst: <b>BCN</b>   |          |
| Percent Moisture                                 | 17     | 1.0     | 1.0    |      | wt%      | 1   | 5/23/2016             | R34494   |
| <b>EPA METHOD 7471: MERCURY</b>                  |        |         |        |      |          |     | Analyst: <b>pmf</b>   |          |
| Mercury  | ND     | 0.00067 | 0.039  |      | mg/Kg-dr | 1   | 5/26/2016 3:51:37 PM  | 25509    |
| <b>EPA METHOD 6010B: SOIL METALS</b>             |        |         |        |      |          |     | Analyst: <b>MED</b>   |          |
| Antimony   | ND     | 2.9     | 6.1    |      | mg/Kg-dr | 2   | 5/31/2016 10:19:25 AM | 25508    |
| Arsenic  | 2.6    | 0.88    | 3.0    | J    | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Barium   | 350    | 0.12    | 0.24   |      | mg/Kg-dr | 2   | 5/31/2016 10:19:25 AM | 25508    |
| Beryllium  | 0.97   | 0.0056  | 0.18   |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Cadmium  | ND     | 0.077   | 0.12   |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Chromium   | 13     | 0.15    | 0.36   |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Cobalt   | 7.8    | 0.046   | 0.36   |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Iron   | 17000  | 120     | 300    |      | mg/Kg-dr | 100 | 5/31/2016 10:17:54 AM | 25508    |
| Lead   | 3.2    | 0.21    | 0.30   |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Manganese  | 520    | 0.11    | 0.24   |      | mg/Kg-dr | 2   | 5/31/2016 10:19:25 AM | 25508    |
| Nickel   | 15     | 0.14    | 0.61   |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Selenium   | ND     | 1.3     | 3.0    |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Silver   | ND     | 0.038   | 0.30   |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Vanadium   | 17     | 0.042   | 3.0    |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| Zinc   | 23     | 0.69    | 3.0    |      | mg/Kg-dr | 1   | 5/26/2016 4:13:12 PM  | 25508    |
| <b>EPA METHOD 8270C: SEMIVOLATILES</b>           |        |         |        |      |          |     | Analyst: <b>DAM</b>   |          |
| Acenaphthene                                     | ND     | 0.10    | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Acenaphthylene                                   | ND     | 0.097   | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Aniline  | ND     | 0.11    | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Anthracene                                       | ND     | 0.079   | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Azobenzene                                       | ND     | 0.15    | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Benz(a)anthracene                                | ND     | 0.10    | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Benzo(a)pyrene                                   | ND     | 0.091   | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Benzo(b)fluoranthene                             | ND     | 0.11    | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Benzo(g,h,i)perylene                             | ND     | 0.11    | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Benzo(k)fluoranthene                             | ND     | 0.11    | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Benzoic acid                                     | ND     | 0.099   | 0.60   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Benzyl alcohol                                   | ND     | 0.094   | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |
| Bis(2-chloroethoxy)methane                       | ND     | 0.13    | 0.24   |      | mg/Kg-dr | 1   | 5/26/2016 8:23:56 PM  | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-22 (8-9')

Project: SWMU 10

Collection Date: 5/18/2016 4:50:00 PM

Lab ID: 1605943-005

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units    | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |      |      |          |    | Analyst: DAM         |          |
| Bis(2-chloroethyl)ether                | ND     | 0.088 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Bis(2-chloroisopropyl)ether            | ND     | 0.11  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Bis(2-ethylhexyl)phthalate             | 0.17   | 0.098 | 0.60 | J    | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 4-Bromophenyl phenyl ether             | ND     | 0.11  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Butyl benzyl phthalate                 | ND     | 0.11  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Carbazole                              | ND     | 0.081 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 4-Chloro-3-methylphenol                | ND     | 0.14  | 0.60 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 4-Chloroaniline                        | ND     | 0.13  | 0.60 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2-Chloronaphthalene                    | ND     | 0.094 | 0.30 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2-Chlorophenol                         | ND     | 0.094 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 4-Chlorophenyl phenyl ether            | ND     | 0.14  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Chrysene                               | ND     | 0.10  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Di-n-butyl phthalate                   | 0.11   | 0.089 | 0.48 | J    | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Di-n-octyl phthalate                   | ND     | 0.10  | 0.48 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Dibenz(a,h)anthracene                  | ND     | 0.097 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Dibenzofuran                           | ND     | 0.12  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 1,2-Dichlorobenzene                    | ND     | 0.092 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 1,3-Dichlorobenzene                    | ND     | 0.093 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 1,4-Dichlorobenzene                    | ND     | 0.10  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 3,3'-Dichlorobenzidine                 | ND     | 0.088 | 0.30 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Diethyl phthalate                      | ND     | 0.12  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Dimethyl phthalate                     | ND     | 0.12  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2,4-Dichlorophenol                     | ND     | 0.11  | 0.48 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2,4-Dimethylphenol                     | ND     | 0.13  | 0.36 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 4,6-Dinitro-2-methylphenol             | ND     | 0.072 | 0.48 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2,4-Dinitrophenol                      | ND     | 0.079 | 0.60 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2,4-Dinitrotoluene                     | ND     | 0.11  | 0.60 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2,6-Dinitrotoluene                     | ND     | 0.13  | 0.60 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Fluoranthene                           | ND     | 0.069 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Fluorene                               | ND     | 0.11  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Hexachlorobenzene                      | ND     | 0.094 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Hexachlorobutadiene                    | ND     | 0.13  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Hexachlorocyclopentadiene              | ND     | 0.14  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Hexachloroethane                       | ND     | 0.10  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Indeno(1,2,3-cd)pyrene                 | ND     | 0.093 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 1-Methylnaphthalene                    | ND     | 0.12  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2-Methylnaphthalene                    | ND     | 0.14  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2-Methylphenol                         | ND     | 0.10  | 0.48 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 3+4-Methylphenol                       | ND     | 0.087 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:23:56 PM | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |   |   |
|--------------------|---|---|
| <b>Qualifiers:</b> | * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
|                    | D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
|                    | H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
|                    | ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
|                    | R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
|                    | S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-22 (8-9')

**Project:** SWMU 10

**Collection Date:** 5/18/2016 4:50:00 PM

**Lab ID:** 1605943-005

**Matrix:** SOIL

**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL      | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|----------|------|-----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |          |      |           |    | Analyst: <b>DAM</b>  |          |
| N-Nitrosodi-n-propylamine              | ND     | 0.12  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| N-Nitrosodiphenylamine                 | ND     | 0.12  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Naphthalene                            | ND     | 0.11  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2-Nitroaniline                         | ND     | 0.13  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 3-Nitroaniline                         | ND     | 0.11  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 4-Nitroaniline                         | ND     | 0.084 | 0.48     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Nitrobenzene                           | ND     | 0.12  | 0.48     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2-Nitrophenol                          | ND     | 0.12  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 4-Nitrophenol                          | ND     | 0.091 | 0.30     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Pentachlorophenol                      | ND     | 0.077 | 0.48     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Phenanthrene                           | ND     | 0.081 | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Phenol                                 | ND     | 0.090 | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Pyrene                                 | ND     | 0.090 | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Pyridine                               | ND     | 0.095 | 0.48     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 1,2,4-Trichlorobenzene                 | ND     | 0.13  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2,4,5-Trichlorophenol                  | ND     | 0.12  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| 2,4,6-Trichlorophenol                  | ND     | 0.099 | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Surr: 2-Fluorophenol                   | 54.0   | 0     | 28.3-102 |      | %Rec      | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Surr: Phenol-d5                        | 55.0   | 0     | 35.7-103 |      | %Rec      | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Surr: 2,4,6-Tribromophenol             | 68.4   | 0     | 35.2-108 |      | %Rec      | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Surr: Nitrobenzene-d5                  | 61.5   |       | 24-118   |      | %Rec      | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Surr: 2-Fluorobiphenyl                 | 67.3   |       | 35.4-111 |      | %Rec      | 1  | 5/26/2016 8:23:56 PM | 25469    |
| Surr: 4-Terphenyl-d14                  | 59.9   |       | 15-91.7  |      | %Rec      | 1  | 5/26/2016 8:23:56 PM | 25469    |
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |          |      |           |    | Analyst: <b>BCN</b>  |          |
| Benzene                                | 9.54   | 2.22  | 2.22     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Toluene                                | 7.02   | 0.268 | 2.22     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Ethylbenzene                           | 1.43   | 0.298 | 2.22     | J    | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Methyl tert-butyl ether (MTBE)         | ND     | 0.373 | 2.22     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,2,4-Trimethylbenzene                 | 1.08   | 0.376 | 2.22     | J    | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,3,5-Trimethylbenzene                 | 0.677  | 0.366 | 2.22     | J    | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,2-Dichloroethane (EDC)               | ND     | 2.22  | 2.22     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,2-Dibromoethane (EDB)                | ND     | 2.22  | 2.22     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Naphthalene                            | ND     | 2.22  | 2.22     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1-Methylnaphthalene                    | ND     | 0.252 | 4.44     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 2-Methylnaphthalene                    | ND     | 0.584 | 4.44     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Acetone                                | 2.93   | 0.657 | 11.1     | J    | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Bromobenzene                           | ND     | 0.226 | 2.22     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Bromodichloromethane                   | ND     | 2.22  | 2.22     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Bromoform                              | ND     | 2.22  | 2.22     |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

**Hall Environmental Analysis Laboratory, Inc.****CLIENT:** Western Refining Company**Client Sample ID:** SWMU 10-22 (8-9')**Project:** SWMU 10**Collection Date:** 5/18/2016 4:50:00 PM**Lab ID:** 1605943-005**Matrix:** SOIL**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |      |      |           |    | Analyst: <b>BCN</b>  |          |
| Bromomethane                           | ND     | 0.400 | 3.33 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 2-Butanone                             | ND     | 0.800 | 11.1 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Carbon disulfide                       | ND     | 0.821 | 11.1 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Carbon tetrachloride                   | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Chlorobenzene                          | ND     | 0.255 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Chloroethane                           | ND     | 0.410 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Chloroform                             | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Chloromethane                          | ND     | 0.562 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 2-Chlorotoluene                        | ND     | 0.378 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 4-Chlorotoluene                        | ND     | 0.369 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| cis-1,2-DCE                            | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| cis-1,3-Dichloropropene                | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,2-Dibromo-3-chloropropane            | ND     | 0.236 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Dibromochloromethane                   | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Dibromomethane                         | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,2-Dichlorobenzene                    | ND     | 0.323 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,3-Dichlorobenzene                    | ND     | 0.410 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,4-Dichlorobenzene                    | ND     | 0.415 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Dichlorodifluoromethane                | ND     | 1.30  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,1-Dichloroethane                     | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,1-Dichloroethene                     | ND     | 0.363 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,2-Dichloropropane                    | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,3-Dichloropropane                    | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 2,2-Dichloropropane                    | ND     | 0.280 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,1-Dichloropropene                    | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Hexachlorobutadiene                    | ND     | 0.506 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 2-Hexanone                             | ND     | 0.553 | 11.1 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Isopropylbenzene                       | ND     | 0.279 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 4-Isopropyltoluene                     | ND     | 0.416 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 4-Methyl-2-pentanone                   | ND     | 4.44  | 11.1 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Methylene chloride                     | ND     | 2.22  | 3.33 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| n-Butylbenzene                         | ND     | 0.540 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| n-Propylbenzene                        | ND     | 0.397 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| sec-Butylbenzene                       | ND     | 0.394 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Styrene                                | ND     | 0.284 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| tert-Butylbenzene                      | ND     | 0.324 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,1,1,2-Tetrachloroethane              | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,1,2,2-Tetrachloroethane              | ND     | 2.22  | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Tetrachloroethene (PCE)                | ND     | 0.292 | 2.22 |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-22 (8-9')

**Project:** SWMU 10

**Collection Date:** 5/18/2016 4:50:00 PM

**Lab ID:** 1605943-005

**Matrix:** SOIL

**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                                    | Result | MDL   | PQL    | Qual | Units     | DF | Date Analyzed        | Batch ID |
|---|--------|-------|--------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b>      |        |       |        |      |           |    | Analyst: <b>BCN</b>  |          |
| trans-1,2-DCE                               | ND     | 0.224 | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| trans-1,3-Dichloropropene                   | ND     | 0.269 | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,2,3-Trichlorobenzene                      | ND     | 0.537 | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,2,4-Trichlorobenzene                      | ND     | 0.673 | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,1,1-Trichloroethane                       | ND     | 2.22  | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,1,2-Trichloroethane                       | ND     | 2.22  | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Trichloroethene (TCE)                       | ND     | 2.22  | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Trichlorofluoromethane                      | ND     | 0.279 | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| 1,2,3-Trichloropropane                      | ND     | 2.22  | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Vinyl chloride                              | ND     | 0.584 | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Xylenes, Total                              | 4.19   | 0.885 | 2.22   |      | µg/Kg-dry | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Surr: 1,2-Dichloroethane-d4                 | 101    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Surr: 4-Bromofluorobenzene                  | 99.3   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Surr: Dibromofluoromethane                  | 90.9   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 6:39:00 PM | 25468    |
| Surr: Toluene-d8                            | 96.4   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 6:39:00 PM | 25468    |
| <b>EPA METHOD 8015D MOD: GASOLINE RANGE</b> |        |       |        |      |           |    | Analyst: <b>DJF</b>  |          |
| Gasoline Range Organics (GRO)               | 0.73   | 0.52  | 3.4    | J    | mg/Kg-dr  | 1  | 5/24/2016 2:30:18 PM | A34454   |
| Surr: BFB                                   | 97.7   | 0     | 70-130 |      | %Rec      | 1  | 5/24/2016 2:30:18 PM | A34454   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |   |   |
|--------------------|---|---|
| <b>Qualifiers:</b> | * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
|                    | D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
|                    | H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
|                    | ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
|                    | R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
|                    | S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |



# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: DUP01

Project: SWMU 10

Collection Date: 5/18/2016

Lab ID: 1605943-006

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses   | Result | MDL     | PQL    | Qual | Units    | DF | Date Analyzed         | Batch ID |
|--|--------|---------|--------|------|----------|----|-----------------------|----------|
| <b>EPA METHOD 8015M/D: DIESEL RANGE ORGANICS</b> |        |         |        |      |          |    | Analyst: <b>KJH</b>   |          |
| Diesel Range Organics (DRO)                      | ND     | 4.3     | 12     |      | mg/Kg-dr | 1  | 5/25/2016 9:22:37 AM  | 25449    |
| Motor Oil Range Organics (MRO)                   | ND     | 60      | 60     |      | mg/Kg-dr | 1  | 5/25/2016 9:22:37 AM  | 25449    |
| Surr: DNOP                                       | 106    | 0       | 70-130 |      | %Rec     | 1  | 5/25/2016 9:22:37 AM  | 25449    |
| <b>PERCENT MOISTURE</b>                          |        |         |        |      |          |    | Analyst: <b>BCN</b>   |          |
| Percent Moisture                                 | 19     | 1.0     | 1.0    |      | wt%      | 1  | 5/23/2016             | R34494   |
| <b>EPA METHOD 7471: MERCURY</b>                  |        |         |        |      |          |    | Analyst: <b>pmf</b>   |          |
| Mercury  | 0.018  | 0.00071 | 0.041  | J    | mg/Kg-dr | 1  | 5/26/2016 3:53:25 PM  | 25509    |
| <b>EPA METHOD 6010B: SOIL METALS</b>             |        |         |        |      |          |    | Analyst: <b>MED</b>   |          |
| Antimony   | ND     | 7.3     | 15     |      | mg/Kg-dr | 5  | 5/31/2016 10:28:43 AM | 25508    |
| Arsenic  | 2.2    | 0.90    | 3.1    | J    | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Barium   | 730    | 0.30    | 0.62   |      | mg/Kg-dr | 5  | 5/31/2016 10:28:43 AM | 25508    |
| Beryllium  | 0.90   | 0.0057  | 0.18   |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Cadmium  | ND     | 0.078   | 0.12   |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Chromium   | 9.8    | 0.15    | 0.37   |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Cobalt   | 4.4    | 0.047   | 0.37   |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Iron   | 14000  | 60      | 150    |      | mg/Kg-dr | 50 | 5/31/2016 10:21:09 AM | 25508    |
| Lead   | 4.6    | 0.22    | 0.31   |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Manganese  | 510    | 0.27    | 0.62   |      | mg/Kg-dr | 5  | 5/31/2016 10:28:43 AM | 25508    |
| Nickel   | 8.0    | 0.14    | 0.62   |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Selenium   | ND     | 1.3     | 3.1    |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Silver   | ND     | 0.039   | 0.31   |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Vanadium   | 19     | 0.043   | 3.1    |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| Zinc   | 24     | 0.70    | 3.1    |      | mg/Kg-dr | 1  | 5/26/2016 4:14:40 PM  | 25508    |
| <b>EPA METHOD 8270C: SEMIVOLATILES</b>           |        |         |        |      |          |    | Analyst: <b>DAM</b>   |          |
| Acenaphthene                                     | ND     | 0.10    | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Acenaphthylene                                   | ND     | 0.099   | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Aniline  | ND     | 0.11    | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Anthracene                                       | ND     | 0.081   | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Azobenzene                                       | ND     | 0.15    | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Benz(a)anthracene                                | ND     | 0.10    | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Benzo(a)pyrene                                   | ND     | 0.092   | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Benzo(b)fluoranthene                             | ND     | 0.11    | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Benzo(g,h,i)perylene                             | ND     | 0.11    | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Benzo(k)fluoranthene                             | ND     | 0.11    | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Benzoic acid                                     | ND     | 0.10    | 0.61   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Benzyl alcohol                                   | ND     | 0.095   | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |
| Bis(2-chloroethoxy)methane                       | ND     | 0.13    | 0.24   |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM  | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

## Hall Environmental Analysis Laboratory, Inc.

CLIENT: Western Refining Company

Client Sample ID: DUP01

Project: SWMU 10

Collection Date: 5/18/2016

Lab ID: 1605943-006

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units    | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |      |      |          |    | Analyst: DAM         |          |
| Bis(2-chloroethyl)ether                | ND     | 0.089 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Bis(2-chloroisopropyl)ether            | ND     | 0.11  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Bis(2-ethylhexyl)phthalate             | 0.16   | 0.099 | 0.61 | J    | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 4-Bromophenyl phenyl ether             | ND     | 0.12  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Butyl benzyl phthalate                 | ND     | 0.11  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Carbazole                              | ND     | 0.082 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 4-Chloro-3-methylphenol                | ND     | 0.14  | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 4-Chloroaniline                        | ND     | 0.13  | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2-Chloronaphthalene                    | ND     | 0.096 | 0.30 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2-Chlorophenol                         | ND     | 0.096 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 4-Chlorophenyl phenyl ether            | ND     | 0.14  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Chrysene                               | ND     | 0.10  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Di-n-butyl phthalate                   | ND     | 0.091 | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Di-n-octyl phthalate                   | ND     | 0.10  | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Dibenz(a,h)anthracene                  | ND     | 0.098 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Dibenzofuran                           | ND     | 0.12  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 1,2-Dichlorobenzene                    | ND     | 0.093 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 1,3-Dichlorobenzene                    | ND     | 0.094 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 1,4-Dichlorobenzene                    | ND     | 0.10  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 3,3'-Dichlorobenzidine                 | ND     | 0.089 | 0.30 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Diethyl phthalate                      | ND     | 0.12  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Dimethyl phthalate                     | ND     | 0.12  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2,4-Dichlorophenol                     | ND     | 0.11  | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2,4-Dimethylphenol                     | ND     | 0.13  | 0.36 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 4,6-Dinitro-2-methylphenol             | ND     | 0.073 | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2,4-Dinitrophenol                      | ND     | 0.080 | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2,4-Dinitrotoluene                     | ND     | 0.11  | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2,6-Dinitrotoluene                     | ND     | 0.13  | 0.61 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Fluoranthene                           | ND     | 0.070 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Fluorene                               | ND     | 0.11  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Hexachlorobenzene                      | ND     | 0.096 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Hexachlorobutadiene                    | ND     | 0.14  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Hexachlorocyclopentadiene              | ND     | 0.14  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Hexachloroethane                       | ND     | 0.10  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Indeno(1,2,3-cd)pyrene                 | ND     | 0.095 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 1-Methylnaphthalene                    | ND     | 0.12  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2-Methylnaphthalene                    | ND     | 0.14  | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2-Methylphenol                         | ND     | 0.10  | 0.49 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 3+4-Methylphenol                       | ND     | 0.088 | 0.24 |      | mg/Kg-dr | 1  | 5/26/2016 8:54:00 PM | 25469    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** DUP01

**Project:** SWMU 10

**Collection Date:** 5/18/2016

**Lab ID:** 1605943-006

**Matrix:** SOIL

**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL      | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|----------|------|-----------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |       |          |      |           |    | Analyst: <b>DAM</b>  |          |
| N-Nitrosodi-n-propylamine              | ND     | 0.12  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| N-Nitrosodiphenylamine                 | ND     | 0.12  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Naphthalene                            | ND     | 0.12  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2-Nitroaniline                         | ND     | 0.13  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 3-Nitroaniline                         | ND     | 0.11  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 4-Nitroaniline                         | ND     | 0.086 | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Nitrobenzene                           | ND     | 0.13  | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2-Nitrophenol                          | ND     | 0.12  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 4-Nitrophenol                          | ND     | 0.092 | 0.30     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Pentachlorophenol                      | ND     | 0.078 | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Phenanthrene                           | ND     | 0.082 | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Phenol                                 | ND     | 0.091 | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Pyrene                                 | ND     | 0.092 | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Pyridine                               | ND     | 0.096 | 0.49     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 1,2,4-Trichlorobenzene                 | ND     | 0.13  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2,4,5-Trichlorophenol                  | ND     | 0.12  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| 2,4,6-Trichlorophenol                  | ND     | 0.10  | 0.24     |      | mg/Kg-dr  | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Surr: 2-Fluorophenol                   | 32.2   | 0     | 28.3-102 |      | %Rec      | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Surr: Phenol-d5                        | 37.7   | 0     | 35.7-103 |      | %Rec      | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Surr: 2,4,6-Tribromophenol             | 61.5   | 0     | 35.2-108 |      | %Rec      | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Surr: Nitrobenzene-d5                  | 36.4   |       | 24-118   |      | %Rec      | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Surr: 2-Fluorobiphenyl                 | 45.6   |       | 35.4-111 |      | %Rec      | 1  | 5/26/2016 8:54:00 PM | 25469    |
| Surr: 4-Terphenyl-d14                  | 65.4   |       | 15-91.7  |      | %Rec      | 1  | 5/26/2016 8:54:00 PM | 25469    |
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |          |      |           |    | Analyst: <b>BCN</b>  |          |
| Benzene                                | ND     | 1.80  | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Toluene                                | 0.918  | 0.217 | 1.80     | J    | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Ethylbenzene                           | ND     | 0.242 | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Methyl tert-butyl ether (MTBE)         | ND     | 0.302 | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,2,4-Trimethylbenzene                 | ND     | 0.305 | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,3,5-Trimethylbenzene                 | ND     | 0.297 | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,2-Dichloroethane (EDC)               | ND     | 1.80  | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,2-Dibromoethane (EDB)                | ND     | 1.80  | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Naphthalene                            | ND     | 1.80  | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1-Methylnaphthalene                    | ND     | 0.205 | 3.60     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 2-Methylnaphthalene                    | ND     | 0.474 | 3.60     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Acetone                                | 8.23   | 0.533 | 9.00     | J    | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Bromobenzene                           | ND     | 0.183 | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Bromodichloromethane                   | ND     | 1.80  | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Bromoform                              | ND     | 1.80  | 1.80     |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

**Hall Environmental Analysis Laboratory, Inc.****CLIENT:** Western Refining Company**Client Sample ID:** DUP01**Project:** SWMU 10**Collection Date:** 5/18/2016**Lab ID:** 1605943-006**Matrix:** SOIL**Received Date:** 5/19/2016 4:30:00 PM

| Analyses                               | Result | MDL   | PQL  | Qual | Units     | DF | Date Analyzed        | Batch ID |
|--|--------|-------|------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b> |        |       |      |      |           |    | Analyst: <b>BCN</b>  |          |
| Bromomethane                           | ND     | 0.324 | 2.70 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 2-Butanone                             | ND     | 0.649 | 9.00 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Carbon disulfide                       | ND     | 0.666 | 9.00 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Carbon tetrachloride                   | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Chlorobenzene                          | ND     | 0.206 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Chloroethane                           | ND     | 0.332 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Chloroform                             | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Chloromethane                          | ND     | 0.456 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 2-Chlorotoluene                        | ND     | 0.306 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 4-Chlorotoluene                        | ND     | 0.299 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| cis-1,2-DCE                            | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| cis-1,3-Dichloropropene                | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,2-Dibromo-3-chloropropane            | ND     | 0.192 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Dibromochloromethane                   | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Dibromomethane                         | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,2-Dichlorobenzene                    | ND     | 0.262 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,3-Dichlorobenzene                    | ND     | 0.332 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,4-Dichlorobenzene                    | ND     | 0.337 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Dichlorodifluoromethane                | ND     | 1.05  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,1-Dichloroethane                     | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,1-Dichloroethene                     | ND     | 0.295 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,2-Dichloropropane                    | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,3-Dichloropropane                    | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 2,2-Dichloropropane                    | ND     | 0.227 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,1-Dichloropropene                    | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Hexachlorobutadiene                    | ND     | 0.410 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 2-Hexanone                             | ND     | 0.448 | 9.00 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Isopropylbenzene                       | ND     | 0.226 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 4-Isopropyltoluene                     | ND     | 0.338 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 4-Methyl-2-pentanone                   | ND     | 3.60  | 9.00 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Methylene chloride                     | ND     | 1.80  | 2.70 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| n-Butylbenzene                         | ND     | 0.438 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| n-Propylbenzene                        | ND     | 0.322 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| sec-Butylbenzene                       | ND     | 0.320 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Styrene                                | ND     | 0.230 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| tert-Butylbenzene                      | ND     | 0.263 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,1,1,2-Tetrachloroethane              | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,1,2,2-Tetrachloroethane              | ND     | 1.80  | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Tetrachloroethene (PCE)                | ND     | 0.236 | 1.80 |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: DUP01

Project: SWMU 10

Collection Date: 5/18/2016

Lab ID: 1605943-006

Matrix: SOIL

Received Date: 5/19/2016 4:30:00 PM

| Analyses                                    | Result | MDL   | PQL    | Qual | Units     | DF | Date Analyzed        | Batch ID |
|---|--------|-------|--------|------|-----------|----|----------------------|----------|
| <b>METHOD 8260B/5035LOW: VOLATILES</b>      |        |       |        |      |           |    | Analyst: <b>BCN</b>  |          |
| trans-1,2-DCE                               | ND     | 0.182 | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| trans-1,3-Dichloropropene                   | ND     | 0.218 | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,2,3-Trichlorobenzene                      | ND     | 0.435 | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,2,4-Trichlorobenzene                      | ND     | 0.546 | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,1,1-Trichloroethane                       | ND     | 1.80  | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,1,2-Trichloroethane                       | ND     | 1.80  | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Trichloroethene (TCE)                       | ND     | 1.80  | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Trichlorofluoromethane                      | ND     | 0.226 | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| 1,2,3-Trichloropropane                      | ND     | 1.80  | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Vinyl chloride                              | ND     | 0.473 | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Xylenes, Total                              | ND     | 0.717 | 1.80   |      | µg/Kg-dry | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Surr: 1,2-Dichloroethane-d4                 | 103    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Surr: 4-Bromofluorobenzene                  | 101    | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Surr: Dibromofluoromethane                  | 92.5   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 7:04:00 PM | 25468    |
| Surr: Toluene-d8                            | 96.7   | 0     | 70-130 |      | %Rec      | 1  | 5/25/2016 7:04:00 PM | 25468    |
| <b>EPA METHOD 8015D MOD: GASOLINE RANGE</b> |        |       |        |      |           |    | Analyst: <b>DJF</b>  |          |
| Gasoline Range Organics (GRO)               | 0.70   | 0.51  | 3.4    | J    | mg/Kg-dr  | 1  | 5/24/2016 2:58:36 PM | A34454   |
| Surr: BFB                                   | 99.9   | 0     | 70-130 |      | %Rec      | 1  | 5/24/2016 2:58:36 PM | A34454   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: MeOH Blank

Project: SWMU 10

Collection Date:

Lab ID: 1605943-007

Matrix: MEOH BLAN

Received Date: 5/19/2016 4:30:00 PM

| Analyses                           | Result | MDL    | PQL   | Qual | Units | DF | Date Analyzed        | Batch ID |
|------------------------------------|--------|--------|-------|------|-------|----|----------------------|----------|
| <b>EPA METHOD 8260B: VOLATILES</b> |        |        |       |      |       |    | Analyst: DJF         |          |
| Benzene                            | ND     | 0.020  | 0.025 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Toluene                            | ND     | 0.0030 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Ethylbenzene                       | ND     | 0.0041 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Methyl tert-butyl ether (MTBE)     | ND     | 0.016  | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,2,4-Trimethylbenzene             | ND     | 0.0037 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,3,5-Trimethylbenzene             | ND     | 0.0036 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,2-Dichloroethane (EDC)           | ND     | 0.013  | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,2-Dibromoethane (EDB)            | ND     | 0.0036 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Naphthalene                        | 0.023  | 0.0078 | 0.10  | J    | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1-Methylnaphthalene                | 0.050  | 0.011  | 0.20  | J    | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 2-Methylnaphthalene                | 0.060  | 0.011  | 0.20  | J    | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Acetone                            | ND     | 0.065  | 0.75  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Bromobenzene                       | ND     | 0.0040 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Bromodichloromethane               | ND     | 0.0029 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Bromoform                          | ND     | 0.0061 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Bromomethane                       | ND     | 0.018  | 0.15  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 2-Butanone                         | ND     | 0.029  | 0.50  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Carbon disulfide                   | ND     | 0.017  | 0.50  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Carbon tetrachloride               | ND     | 0.0033 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Chlorobenzene                      | ND     | 0.0041 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Chloroethane                       | ND     | 0.010  | 0.10  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Chloroform                         | ND     | 0.0038 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Chloromethane                      | ND     | 0.0044 | 0.15  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 2-Chlorotoluene                    | ND     | 0.0037 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 4-Chlorotoluene                    | ND     | 0.0044 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| cis-1,2-DCE                        | ND     | 0.0029 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| cis-1,3-Dichloropropene            | ND     | 0.0046 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,2-Dibromo-3-chloropropane        | ND     | 0.015  | 0.10  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Dibromochloromethane               | ND     | 0.0045 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Dibromomethane                     | ND     | 0.0043 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,2-Dichlorobenzene                | ND     | 0.0044 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,3-Dichlorobenzene                | ND     | 0.0041 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,4-Dichlorobenzene                | ND     | 0.0062 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Dichlorodifluoromethane            | ND     | 0.015  | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,1-Dichloroethane                 | ND     | 0.0027 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,1-Dichloroethene                 | ND     | 0.016  | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,2-Dichloropropane                | ND     | 0.0042 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,3-Dichloropropane                | ND     | 0.0057 | 0.050 |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 2,2-Dichloropropane                | ND     | 0.0029 | 0.10  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605943

Date Reported: 6/28/2016

CLIENT: Western Refining Company

Client Sample ID: MeOH Blank

Project: SWMU 10

Collection Date:

Lab ID: 1605943-007

Matrix: MEOH BLAN

Received Date: 5/19/2016 4:30:00 PM

| Analyses                           | Result | MDL    | PQL    | Qual | Units | DF | Date Analyzed        | Batch ID |
|------------------------------------|--------|--------|--------|------|-------|----|----------------------|----------|
| <b>EPA METHOD 8260B: VOLATILES</b> |        |        |        |      |       |    | Analyst: DJF         |          |
| 1,1-Dichloropropene                | ND     | 0.0040 | 0.10   |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Hexachlorobutadiene                | ND     | 0.0061 | 0.10   |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 2-Hexanone                         | ND     | 0.027  | 0.50   |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Isopropylbenzene                   | ND     | 0.0043 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 4-Isopropyltoluene                 | ND     | 0.0045 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 4-Methyl-2-pentanone               | ND     | 0.015  | 0.50   |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Methylene chloride                 | 0.016  | 0.014  | 0.15   | J    | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| n-Butylbenzene                     | ND     | 0.0044 | 0.15   |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| n-Propylbenzene                    | ND     | 0.0038 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| sec-Butylbenzene                   | ND     | 0.0069 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Styrene                            | ND     | 0.0045 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| tert-Butylbenzene                  | ND     | 0.0041 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,1,1,2-Tetrachloroethane          | ND     | 0.0048 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,1,2,2-Tetrachloroethane          | ND     | 0.0081 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Tetrachloroethene (PCE)            | ND     | 0.0041 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| trans-1,2-DCE                      | ND     | 0.014  | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| trans-1,3-Dichloropropene          | ND     | 0.0073 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,2,3-Trichlorobenzene             | ND     | 0.0075 | 0.10   |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,2,4-Trichlorobenzene             | ND     | 0.0053 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,1,1-Trichloroethane              | ND     | 0.0031 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,1,2-Trichloroethane              | ND     | 0.0059 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Trichloroethene (TCE)              | ND     | 0.0054 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Trichlorofluoromethane             | ND     | 0.0037 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| 1,2,3-Trichloropropane             | ND     | 0.0086 | 0.10   |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Vinyl chloride                     | ND     | 0.0041 | 0.050  |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Xylenes, Total                     | ND     | 0.0095 | 0.10   |      | mg/Kg | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Surr: Dibromofluoromethane         | 86.8   |        | 70-130 |      | %Rec  | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Surr: 1,2-Dichloroethane-d4        | 78.8   |        | 70-130 |      | %Rec  | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Surr: Toluene-d8                   | 99.5   |        | 70-130 |      | %Rec  | 1  | 5/24/2016 3:55:15 PM | C34454   |
| Surr: 4-Bromofluorobenzene         | 104    |        | 70-130 |      | %Rec  | 1  | 5/24/2016 3:55:15 PM | C34454   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |   |   |
|--------------------|---|---|
| <b>Qualifiers:</b> | * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
|                    | D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
|                    | H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
|                    | ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
|                    | R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
|                    | S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |



## Wet Chemistry by Method 3060A/7196A

| Analyte             | Result | Qualifier | MDL   | RDL  | Dilution | Analysis date / time | Batch    |
|---------------------|--------|-----------|-------|------|----------|----------------------|----------|
| Chromium,Hexavalent | U      |           | 0.640 | 2.00 | 1        | 05/26/2016 19:13     | WG874409 |

## Wet Chemistry by Method 9012B

| Analyte | Result | Qualifier | MDL    | RDL   | Dilution | Analysis date / time | Batch    |
|---------|--------|-----------|--------|-------|----------|----------------------|----------|
| Cyanide | 2.20   |           | 0.0390 | 0.250 | 1        | 05/31/2016 10:25     | WG875385 |

Tc

Ss

Cn

Sr

Qc

Gl

Al

Sc





## Wet Chemistry by Method 3060A/7196A

| Analyte              | Result | Qualifier | MDL   | RDL   | Dilution | Analysis         | Batch    |
|----------------------|--------|-----------|-------|-------|----------|------------------|----------|
|                      | mg/kg  |           | mg/kg | mg/kg |          | date / time      |          |
| Chromium, Hexavalent | U      |           | 0.640 | 2.00  | 1        | 05/26/2016 19:13 | WG874403 |

## Wet Chemistry by Method 9012B

| Analyte | Result | Qualifier | MDL    | RDL   | Dilution | Analysis         | Batch    |
|---------|--------|-----------|--------|-------|----------|------------------|----------|
|         | mg/kg  |           | mg/kg  | mg/kg |          | date / time      |          |
| Cyanide | 0.0485 | J         | 0.0390 | 0.250 | 1        | 05/31/2016 10:27 | WG875385 |

2  
Tc3  
Ss4  
Cn5  
Sr6  
Qc7  
Gl8  
Al9  
Sc



## Wet Chemistry by Method 3060A/7196A

| Analyte              | Result<br>mg/kg | Qualifier | MDL<br>mg/kg | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch           |
|----------------------|-----------------|-----------|--------------|--------------|----------|-------------------------|-----------------|
| Chromium, Hexavalent | U               |           | 0.640        | 2.00         | 1        | 05/26/2016 19:15        | <u>WG874409</u> |

## Wet Chemistry by Method 9012B

| Analyte | Result<br>mg/kg | Qualifier | MDL<br>mg/kg | RDL<br>mg/kg | Dilution | Analysis<br>date / time | Batch           |
|---------|-----------------|-----------|--------------|--------------|----------|-------------------------|-----------------|
| Cyanide | 0.0438          | J         | 0.0390       | 0.250        | 1        | 06/02/2016 09:27        | <u>WG876669</u> |

2 Tc

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



## Wet Chemistry by Method 3060A/7196A

|                      | Result | Qualifier | MDL   | RDL   | Dilution | Analysis<br>date / time | Batch    |
|----------------------|--------|-----------|-------|-------|----------|-------------------------|----------|
| Analyte              | mg/kg  |           | mg/kg | mg/kg |          |                         |          |
| Chromium, Hexavalent | U      |           | 0.640 | 2.00  | 1        | 05/26/2016 19:16        | WG874409 |

## Wet Chemistry by Method 9012B

|         | Result | Qualifier | MDL    | RDL   | Dilution | Analysis<br>date / time | Batch    |
|---------|--------|-----------|--------|-------|----------|-------------------------|----------|
| Analyte | mg/kg  |           | mg/kg  | mg/kg |          |                         |          |
| Cyanide | 0.0403 | J         | 0.0390 | 0.250 | 1        | 06/02/2016 09:28        | WG876669 |

<sup>2</sup> Tc<sup>3</sup> Ss<sup>4</sup> Cn<sup>5</sup> Sr<sup>6</sup> Qc<sup>7</sup> Gl<sup>8</sup> Al<sup>9</sup> Sc



## Wet Chemistry by Method 3060A/7196A

| Analyte              | Result | Qualifier | MDL   | RDL  | Dilution | Analysis date / time | Batch    |
|----------------------|--------|-----------|-------|------|----------|----------------------|----------|
| Chromium, Hexavalent | U      |           | 0.640 | 2.00 | 1        | 05/26/2016 19:16     | WG874403 |

## Wet Chemistry by Method 9012B

| Analyte | Result | Qualifier | MDL    | RDL   | Dilution | Analysis date / time | Batch    |
|---------|--------|-----------|--------|-------|----------|----------------------|----------|
| Cyanide | U      |           | 0.0390 | 0.250 | 1        | 06/02/2016 09:29     | WG876669 |

<sup>2</sup> Tc<sup>3</sup> Ss<sup>4</sup> Cn<sup>5</sup> Sr<sup>6</sup> Qc<sup>7</sup> Gl

Ai

<sup>9</sup> Sc

1605943-006C DUP01

Collected date/time: 05/18/16 00:00

## SAMPLE RESULTS - 06

L837333

ONE LAB. NATIONWIDE.



Wet Chemistry by Method 3060A/7196A

| Analyte              | Result | Qualifier | MDL   | RDL  | Dilution | Analysis<br>date / time | Batch           |
|----------------------|--------|-----------|-------|------|----------|-------------------------|-----------------|
| Chromium, Hexavalent | U      |           | 0.640 | 2.00 | 1        | 05/26/2016 19:17        | <u>WG874409</u> |

Wet Chemistry by Method 9012B

| Analyte | Result | Qualifier | MDL    | RDL   | Dilution | Analysis<br>date / time | Batch           |
|---------|--------|-----------|--------|-------|----------|-------------------------|-----------------|
| Cyanide | U      |           | 0.0390 | 0.250 | 1        | 06/02/2016 09:30        | <u>WG876669</u> |

<sup>2</sup> Tc<sup>3</sup> Ss<sup>4</sup> Cn<sup>5</sup> Sr<sup>6</sup> Qc<sup>7</sup> Gl

Al

<sup>9</sup> Sc

## Method Blank (MB)

| (MB) R3140072-1 05/26/16 18:42 |                    |
|--------------------------------|--------------------|
| Analyte                        | MB Result<br>mg/kg |
| Chromium Hexavalent            | U                  |
|                                | MB MDL<br>mg/kg    |
|                                | 0.640              |
|                                | MB RDL<br>mg/kg    |
|                                | 2.00               |

L837120-10 Original Sample (OS) • Duplicate (DUP)

| (OS) L837120-10 05/26/16 18:49 • (DUP) R3140072-4 05/26/16 18:49 |                       |                  |          |                |
|--|-----------------------|------------------|----------|----------------|
| Analyte  | Original Result (dry) | DUP Result (dry) | Dilution | DUP RPD Limits |
|  | mg/kg                 | mg/kg            |          | %              |
| Chromium, Hexavalent   | U                     | ND               | 1        | 0.000          |
|  |                       |                  |          | 20             |

## L837333-03 Original Sample (OS) • Duplicate (DU2)

| (OS) L837333-03 05/26/16 19:15 • (DUP) R3140072-8 05/26/16 19:15 |                          |                     |          |              |                     |
|--|--------------------------|---------------------|----------|--------------|---------------------|
| Analyte  | Original Result<br>mg/kg | DUP Result<br>mg/kg | Dilution | DUP RPD<br>% | DUP RPD Limits<br>% |
| Chromium, Hexavalent   | U                        | ND                  | 1        | 0.000        | 20                  |

## Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCDSD)

| (LCS) R3140072-2 05/26/16 18:43 • (LCS) R3140072-3 05/26/16 18:44 |              |            |            |          |          |             |            |
|---|--------------|------------|------------|----------|----------|-------------|------------|
| Analyte   | Spike Amount | LCS Result | LCS Result | LCS Rec. | LCS Rec. | Rec. Limits | RPD Limits |
|   | mg/kg        | mg/kg      | mg/kg      | %        | %        | %           | %          |
| Chromium Hexavalent   | 56.9         | 60.0       | 62.8       | 105      | 110      | 80.0-120    | 5.00       |
|   |              |            |            |          |          |             | 20         |

...837120-10 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

[illegible]

ACCOUNT: Hall Environmental Analysis Laboratory

PROJECT:

SDG:  
L837333

DATE/TIME: 06/02/16 15:08



Method Blank (MB)

|                                |           |              |        |        |
|--------------------------------|-----------|--------------|--------|--------|
| (MB) R3140567-4 05/31/16 10:11 |           |              |        |        |
| Analyte                        | MB Result | MB Qualifier | MB MDL | MB RDL |
|                                | mg/kg     | mg/kg        | mg/kg  | mg/kg  |
| Cyanide                        | U         | 0.0390       | 0.250  |        |

• L837333-01 Original Sample (OS) • Duplicate (DUP)

(OS) L837333-01 05/31/16 10:25 • (DUP) R3140567-7 05/31/16 10:26

| Analyte | Original Result | DUP Result | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------|-----------------|------------|---------|---------------|----------------|
|         | mg/kg           | mg/kg      | %       | %             | %              |
| Cyanide | 2.20            | 2.52       | 1       | 13.0          | 20             |

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3140567-5 05/31/16 10:12 • (LCSD) R3140567-6 05/31/16 10:13

| Analyte | Spike Amount | LCS Result | LCSD Result | LCS Rec. | LCSD Rec. | Rec. Limits | LCS Qualifier | LCSD Qualifier | RPD | RPD Limits |
|---------|--------------|------------|-------------|----------|-----------|-------------|---------------|----------------|-----|------------|
|         | mg/kg        | mg/kg      | mg/kg       | %        | %         | %           | %             | %              | %   | %          |
| Cyanide | 48.4         | 66.4       | 59.4        | 137      | 123       | 50.0-150    |               | 11.0           |     | 20         |

|   |    |
|---|----|
| 1 | 1  |
| 3 | SS |
| 4 | Cn |
| 5 | S- |
| 6 | Qc |
| 7 | Gl |
| 8 | AN |
| 9 | SC |

WG876669

Wet Chemistry by Method 9012B

QUALITY CONTROL SUMMARY

L837333-03.04.05.06

ONE LAB. NATIONWIDE



Method Blank (MB)

(MB) R3141092-4 06/02/16 09:22

|         | MB Result | MB Qualifier | MB MDL | MB RDL |
|---------|-----------|--------------|--------|--------|
| Analyte | mg/kg     | mg/kg        | mg/kg  | mg/kg  |
| Cyanide | U         | 0.0390       | 0.250  |        |

L837358-05 Original Sample (OS) • Duplicate (DUP)

(OS) L837358-05 06/02/16 09:36 • (DUP) R3141092-7 06/02/16 09:37

|         | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------|-----------------|------------|----------|---------|---------------|----------------|
| Analyte | mg/kg           | mg/kg      | %        | %       |               | %              |
| Cyanide | 0.104           | ND         | 1        | 47.0    | 3.21          | 20             |

L837397-01 Original Sample (OS) • Duplicate (DUP)

(OS) L837397-01 06/02/16 09:58 • (DUP) R3141092-8 06/02/16 09:59

|         | Original Result | DUP Result | Dilution | DUP RPD | DUP Qualifier | DUP RPD Limits |
|---------|-----------------|------------|----------|---------|---------------|----------------|
| Analyte | mg/kg           | mg/kg      | %        | %       |               | %              |
| Cyanide | ND              | ND         | 10       | 0.000   |               | 20             |

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3141092-5 06/02/16 09:23 • (LCSD) R3141092-6 06/02/16 09:24

|         | Spike Amount | LCS Result | LCSD Result | LCS Rec. | LCSD Rec. | Rec. Limits | LCS Qualifier | LCSD Qualifier | RPD  | RPD Limits |
|---------|--------------|------------|-------------|----------|-----------|-------------|---------------|----------------|------|------------|
| Analyte | mg/kg        | mg/kg      | mg/kg       | %        | %         | %           |               |                | %    | %          |
| Cyanide | 48.4         | 49.0       | 43.2        | 101      | 89.0      | 50.0-150    |               |                | 13.0 | 20         |

L837397-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L837397-03 06/02/16 10:01 • (MS) R3141092-9 06/02/16 10:02 • (MSD) R3141092-10 06/02/16 10:03

|         | Spike Amount (dry) | Original Result (dry) | MS Result (dry) | MSD Result (dry) | MS Rec. | MSD Rec. | Dilution | Rec. Limits | MS Qualifier | MSD Qualifier | RPD  | RPD Limits |
|---------|--------------------|-----------------------|-----------------|------------------|---------|----------|----------|-------------|--------------|---------------|------|------------|
| Analyte | mg/kg              | mg/kg                 | mg/kg           | mg/kg            | %       | %        |          | %           |              |               | %    | %          |
| Cyanide | 3.78               | 1.10                  | 3.76            | 4.23             | 70.0    | 83.0     | 1        | 80.0-120    | 3.6          |               | 12.0 | 20         |

ACCOUNT:  
Hall Environmental Analysis Laboratory

PROJECT:

SDG:  
L837333

DATE/TIME:  
06/02/16 15:08

|   |   |    |    |    |    |    |    |    |
|---|---|----|----|----|----|----|----|----|
| 1 | 2 | 3  | 4  | 5  | 6  | 7  | 8  | 9  |
|   |   | SS | Cn | Sc | Qc | Gl | Al | Sc |





## Abbreviations and Definitions

|                 |   |
|-----------------|---|
| SDG             | Sample Delivery Group.  |
| MDL             | Method Detection Limit.   |
| RDL             | Reported Detection Limit.   |
| ND              | Not detected at the Reporting Limit (or MDL where applicable).  |
| U               | Not detected at the Reporting Limit (or MDL where applicable).  |
| RPD             | Relative Percent Difference.  |
| (dry)           | Results are reported based on the dry weight of the sample. [this will only be present on a dry report basis for soils].  |
| Original Sample | The non-spiked sample in the prep batch used to determine the Relative Percent Difference (RPD) from a quality control sample. The Original Sample may not be included within the reported SDG. |
| Rec.            | Recovery.   |

| Qualifier | Description   |
|-----------|---|
| J         | The identification of the analyte is acceptable; the reported value is an estimate.                   |
| J6        | The sample matrix interfered with the ability to make any accurate determination; spike value is low. |
| P1        | RPD value not applicable for sample concentrations less than 5 times the reporting limit.             |

<sup>2</sup> Tc<sup>3</sup> Ss<sup>4</sup> Cn<sup>5</sup> Sr<sup>6</sup> Qc<sup>7</sup> Gl<sup>8</sup> Al<sup>9</sup> Sc

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                             |           |     |                          |             |   |          |              |      |          |      |
|-----------------------------|-----------|-----|--------------------------|-------------|---|----------|--------------|------|----------|------|
| Sample ID                   | LCS-25449 |     | SampType: LCS            |             | TestCode: EPA Method 8015M/D: Diesel Range Organics |          |              |      |          |      |
| Client ID:                  | LCSS      |     | Batch ID: 25449          |             | RunNo: 34433  |          |              |      |          |      |
| Prep Date:                  | 5/23/2016 |     | Analysis Date: 5/24/2016 |             | SeqNo: 1061812                                      |          | Units: mg/Kg |      |          |      |
| Analyte                     | Result    | PQL | SPK value                | SPK Ref Val | %REC  | LowLimit | HighLimit    | %RPD | RPDLimit | Qual |
| Diesel Range Organics (DRO) | 54        | 10  | 50.00                    | 0           | 109   | 62.6     | 124          |      |          |      |
| Surr: DNOP                  | 5.5       |     | 5.000                    |             | 109   | 70       | 130          |      |          |      |

|                                |           |     |                          |             |   |          |              |      |          |      |
|--------------------------------|-----------|-----|--------------------------|-------------|---|----------|--------------|------|----------|------|
| Sample ID                      | MB-25449  |     | SampType: MBLK           |             | TestCode: EPA Method 8015M/D: Diesel Range Organics |          |              |      |          |      |
| Client ID:                     | PBS       |     | Batch ID: 25449          |             | RunNo: 34433  |          |              |      |          |      |
| Prep Date:                     | 5/23/2016 |     | Analysis Date: 5/24/2016 |             | SeqNo: 1061813                                      |          | Units: mg/Kg |      |          |      |
| Analyte                        | Result    | PQL | SPK value                | SPK Ref Val | %REC  | LowLimit | HighLimit    | %RPD | RPDLimit | Qual |
| Diesel Range Organics (DRO)    | ND        | 10  |                          |             |   |          |              |      |          |      |
| Motor Oil Range Organics (MRO) | ND        | 50  |                          |             |   |          |              |      |          |      |
| Surr: DNOP                     | 11        |     | 10.00                    |             | 113   | 70       | 130          |      |          |      |

### Qualifiers:

|   |   |
|---|---|
| * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
| D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
| H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
| ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
| R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
| S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                                |        |                |           |             |                             |          |           |      |          |      |
|--------------------------------|--------|----------------|-----------|-------------|-----------------------------|----------|-----------|------|----------|------|
| Sample ID                      | rb     | SampType:      | MBLK      | TestCode:   | EPA Method 8260B: Volatiles |          |           |      |          |      |
| Client ID:                     | PBS    | Batch ID:      | C34454    | RunNo:      | 34454                       |          |           |      |          |      |
| Prep Date:                     |        | Analysis Date: | 5/24/2016 | SeqNo:      | 1062796                     | Units:   | mg/Kg     |      |          |      |
| Analyte                        | Result | PQL            | SPK value | SPK Ref Val | %REC                        | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
| Benzene                        | ND     | 0.025          |           |             |                             |          |           |      |          |      |
| Toluene                        | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Ethylbenzene                   | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Methyl tert-butyl ether (MTBE) | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,2,4-Trimethylbenzene         | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,3,5-Trimethylbenzene         | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,2-Dichloroethane (EDC)       | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,2-Dibromoethane (EDB)        | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Naphthalene                    | ND     | 0.10           |           |             |                             |          |           |      |          |      |
| 1-Methylnaphthalene            | ND     | 0.20           |           |             |                             |          |           |      |          |      |
| 2-Methylnaphthalene            | ND     | 0.20           |           |             |                             |          |           |      |          |      |
| Acetone                        | ND     | 0.75           |           |             |                             |          |           |      |          |      |
| Bromobenzene                   | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Bromodichloromethane           | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Bromoform                      | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Bromomethane                   | ND     | 0.15           |           |             |                             |          |           |      |          |      |
| 2-Butanone                     | ND     | 0.50           |           |             |                             |          |           |      |          |      |
| Carbon disulfide               | ND     | 0.50           |           |             |                             |          |           |      |          |      |
| Carbon tetrachloride           | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Chlorobenzene                  | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Chloroethane                   | ND     | 0.10           |           |             |                             |          |           |      |          |      |
| Chloroform                     | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Chloromethane                  | ND     | 0.15           |           |             |                             |          |           |      |          |      |
| 2-Chlorotoluene                | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 4-Chlorotoluene                | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| cis-1,2-DCE                    | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| cis-1,3-Dichloropropene        | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,2-Dibromo-3-chloropropane    | ND     | 0.10           |           |             |                             |          |           |      |          |      |
| Dibromochloromethane           | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Dibromomethane                 | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,2-Dichlorobenzene            | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,3-Dichlorobenzene            | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,4-Dichlorobenzene            | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| Dichlorodifluoromethane        | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,1-Dichloroethane             | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,1-Dichloroethene             | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,2-Dichloropropane            | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 1,3-Dichloropropane            | ND     | 0.050          |           |             |                             |          |           |      |          |      |
| 2,2-Dichloropropane            | ND     | 0.10           |           |             |                             |          |           |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

| Sample ID <b>rb</b>         | SampType: <b>MBLK</b>           |       |           | TestCode: <b>EPA Method 8260B: Volatiles</b> |      |                     |           |      |          |      |
|-----------------------------|---------------------------------|-------|-----------|--|------|---------------------|-----------|------|----------|------|
| Client ID: <b>PBS</b>       | Batch ID: <b>C34454</b>         |       |           | RunNo: <b>34454</b>                          |      |                     |           |      |          |      |
| Prep Date:                  | Analysis Date: <b>5/24/2016</b> |       |           | SeqNo: <b>1062796</b>                        |      | Units: <b>mg/Kg</b> |           |      |          |      |
| Analyte                     | Result                          | PQL   | SPK value | SPK Ref Val                                  | %REC | LowLimit            | HighLimit | %RPD | RPDLimit | Qual |
| 1,1-Dichloropropene         | ND                              | 0.10  |           |  |      |                     |           |      |          |      |
| Hexachlorobutadiene         | ND                              | 0.10  |           |  |      |                     |           |      |          |      |
| 2-Hexanone                  | ND                              | 0.50  |           |  |      |                     |           |      |          |      |
| Isopropylbenzene            | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| 4-Isopropyltoluene          | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| 4-Methyl-2-pentanone        | ND                              | 0.50  |           |  |      |                     |           |      |          |      |
| Methylene chloride          | ND                              | 0.15  |           |  |      |                     |           |      |          |      |
| n-Butylbenzene              | ND                              | 0.15  |           |  |      |                     |           |      |          |      |
| n-Propylbenzene             | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| sec-Butylbenzene            | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| Styrene                     | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| tert-Butylbenzene           | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| 1,1,1,2-Tetrachloroethane   | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| 1,1,2,2-Tetrachloroethane   | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| Tetrachloroethene (PCE)     | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| trans-1,2-DCE               | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| trans-1,3-Dichloropropene   | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| 1,2,3-Trichlorobenzene      | ND                              | 0.10  |           |  |      |                     |           |      |          |      |
| 1,2,4-Trichlorobenzene      | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| 1,1,1-Trichloroethane       | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| 1,1,2-Trichloroethane       | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| Trichloroethene (TCE)       | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| Trichlorofluoromethane      | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| 1,2,3-Trichloropropane      | ND                              | 0.10  |           |  |      |                     |           |      |          |      |
| Vinyl chloride              | ND                              | 0.050 |           |  |      |                     |           |      |          |      |
| Xylenes, Total              | ND                              | 0.10  |           |  |      |                     |           |      |          |      |
| Surr: Dibromofluoromethane  | 0.50                            |       | 0.5000    |  | 99.9 | 70                  | 130       |      |          |      |
| Surr: 1,2-Dichloroethane-d4 | 0.45                            |       | 0.5000    |  | 89.7 | 70                  | 130       |      |          |      |
| Surr: Toluene-d8            | 0.49                            |       | 0.5000    |  | 98.8 | 70                  | 130       |      |          |      |
| Surr: 4-Bromofluorobenzene  | 0.54                            |       | 0.5000    |  | 107  | 70                  | 130       |      |          |      |

| Sample ID <b>100ng lcs</b> | SampType: <b>LCS</b>            |       |           | TestCode: <b>EPA Method 8260B: Volatiles</b> |      |                     |           |      |          |      |
|----------------------------|---------------------------------|-------|-----------|--|------|---------------------|-----------|------|----------|------|
| Client ID: <b>LCSS</b>     | Batch ID: <b>C34454</b>         |       |           | RunNo: <b>34454</b>                          |      |                     |           |      |          |      |
| Prep Date:                 | Analysis Date: <b>5/24/2016</b> |       |           | SeqNo: <b>1062797</b>                        |      | Units: <b>mg/Kg</b> |           |      |          |      |
| Analyte                    | Result                          | PQL   | SPK value | SPK Ref Val                                  | %REC | LowLimit            | HighLimit | %RPD | RPDLimit | Qual |
| Benzene                    | 1.0                             | 0.025 | 1.000     | 0  | 101  | 70                  | 130       |      |          |      |
| Toluene                    | 0.97                            | 0.050 | 1.000     | 0  | 97.4 | 70                  | 130       |      |          |      |
| Chlorobenzene              | 0.95                            | 0.050 | 1.000     | 0  | 95.4 | 70                  | 130       |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

R RPD outside accepted recovery limits

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

P Sample pH Not In Range

RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                             |           |                |           |             |                             |          |           |      |          |      |
|-----------------------------|-----------|----------------|-----------|-------------|-----------------------------|----------|-----------|------|----------|------|
| Sample ID                   | 100ng lcs | SampType:      | LCS       | TestCode:   | EPA Method 8260B: Volatiles |          |           |      |          |      |
| Client ID:                  | LCSS      | Batch ID:      | C34454    | RunNo:      | 34454                       |          |           |      |          |      |
| Prep Date:                  |           | Analysis Date: | 5/24/2016 | SeqNo:      | 1062797                     | Units:   | mg/Kg     |      |          |      |
| Analyte                     | Result    | PQL            | SPK value | SPK Ref Val | %REC                        | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
| 1,1-Dichloroethene          | 1.1       | 0.050          | 1.000     | 0           | 106                         | 70       | 130       |      |          |      |
| Trichloroethene (TCE)       | 0.98      | 0.050          | 1.000     | 0           | 98.2                        | 70       | 130       |      |          |      |
| Surr: Dibromofluoromethane  | 0.51      |                | 0.5000    |             | 102                         | 70       | 130       |      |          |      |
| Surr: 1,2-Dichloroethane-d4 | 0.45      |                | 0.5000    |             | 90.0                        | 70       | 130       |      |          |      |
| Surr: Toluene-d8            | 0.49      |                | 0.5000    |             | 98.0                        | 70       | 130       |      |          |      |
| Surr: 4-Bromofluorobenzene  | 0.54      |                | 0.5000    |             | 109                         | 70       | 130       |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                             |           |      |           |                |           |          |           |                                 |          |        |       |
|-----------------------------|-----------|------|-----------|----------------|-----------|----------|-----------|---------------------------------|----------|--------|-------|
| Sample ID                   | LCS-25468 |      |           | SampType:      | LCS       |          | TestCode: | Method 8260B/5035LOW: VOLATILES |          |        |       |
| Client ID:                  | LCSS      |      |           | Batch ID:      | 25468     |          | RunNo:    | 34484                           |          |        |       |
| Prep Date:                  | 5/24/2016 |      |           | Analysis Date: | 5/25/2016 |          | SeqNo:    | 1063539                         |          | Units: | µg/Kg |
| Analyte                     | Result    | PQL  | SPK value | SPK Ref Val    | %REC      | LowLimit | HighLimit | %RPD                            | RPDLimit | Qual   |       |
| Benzene                     | 10.2      | 2.00 | 10.00     | 0              | 102       | 70       | 130       |                                 |          |        |       |
| Toluene                     | 11.1      | 2.00 | 10.00     | 0              | 111       | 70       | 130       |                                 |          |        |       |
| Chlorobenzene               | 11.1      | 2.00 | 10.00     | 0              | 111       | 70       | 130       |                                 |          |        |       |
| 1,1-Dichloroethene          | 10.3      | 2.00 | 10.00     | 0              | 103       | 68       | 129       |                                 |          |        |       |
| Trichloroethene (TCE)       | 10.2      | 2.00 | 10.00     | 0              | 102       | 70       | 130       |                                 |          |        |       |
| Surr: 1,2-Dichloroethane-d4 | 9.00      |      | 10.00     |                | 90.0      | 70       | 130       |                                 |          |        |       |
| Surr: 4-Bromofluorobenzene  | 10.0      |      | 10.00     |                | 100       | 70       | 130       |                                 |          |        |       |
| Surr: Dibromofluoromethane  | 8.76      |      | 10.00     |                | 87.6      | 70       | 130       |                                 |          |        |       |
| Surr: Toluene-d8            | 9.90      |      | 10.00     |                | 99.0      | 70       | 130       |                                 |          |        |       |

|                             |            |      |                          |             |   |          |              |       |          |      |
|-----------------------------|------------|------|--------------------------|-------------|---|----------|--------------|-------|----------|------|
| Sample ID                   | LCSD-25468 |      | SampType: LCS            |             | TestCode: Method 8260B/5035LOW: VOLATILES |          |              |       |          |      |
| Client ID:                  | LCSS       |      | Batch ID: 25468          |             | RunNo: 34484                              |          |              |       |          |      |
| Prep Date:                  | 5/24/2016  |      | Analysis Date: 5/25/2016 |             | SeqNo: 1063540                            |          | Units: µg/Kg |       |          |      |
| Analyte                     | Result     | PQL  | SPK value                | SPK Ref Val | %REC                                      | LowLimit | HighLimit    | %RPD  | RPDLimit | Qual |
| Benzene                     | 10.3       | 2.00 | 10.00                    | 0           | 103                                       | 70       | 130          | 1.08  | 0        |      |
| Toluene                     | 11.2       | 2.00 | 10.00                    | 0           | 112                                       | 70       | 130          | 0.629 | 0        |      |
| Chlorobenzene               | 11.3       | 2.00 | 10.00                    | 0           | 113                                       | 70       | 130          | 1.96  | 0        |      |
| 1,1-Dichloroethene          | 10.4       | 2.00 | 10.00                    | 0           | 104                                       | 68       | 129          | 1.16  | 0        |      |
| Trichloroethene (TCE)       | 10.2       | 2.00 | 10.00                    | 0           | 102                                       | 70       | 130          | 0.295 | 0        |      |
| Surr: 1,2-Dichloroethane-d4 | 8.93       |      | 10.00                    |             | 89.3                                      | 70       | 130          | 0     | 0        |      |
| Surr: 4-Bromofluorobenzene  | 10.1       |      | 10.00                    |             | 101                                       | 70       | 130          | 0     | 0        |      |
| Surr: Dibromofluoromethane  | 8.72       |      | 10.00                    |             | 87.2                                      | 70       | 130          | 0     | 0        |      |
| Surr: Toluene-d8            | 9.86       |      | 10.00                    |             | 98.6                                      | 70       | 130          | 0     | 0        |      |

|                                |           |      |                          |             |   |          |              |      |          |      |
|--------------------------------|-----------|------|--------------------------|-------------|---|----------|--------------|------|----------|------|
| Sample ID                      | MB-25468  |      | SampType: MBLK           |             | TestCode: Method 8260B/5035LOW: VOLATILES |          |              |      |          |      |
| Client ID:                     | PBS       |      | Batch ID: 25468          |             | RunNo: 34484                              |          |              |      |          |      |
| Prep Date:                     | 5/24/2016 |      | Analysis Date: 5/25/2016 |             | SeqNo: 1063541                            |          | Units: µg/Kg |      |          |      |
| Analyte                        | Result    | PQL  | SPK value                | SPK Ref Val | %REC                                      | LowLimit | HighLimit    | %RPD | RPDLimit | Qual |
| Benzene                        | ND        | 2.00 |                          |             |   |          |              |      |          |      |
| Toluene                        | ND        | 2.00 |                          |             |   |          |              |      |          |      |
| Ethylbenzene                   | ND        | 2.00 |                          |             |   |          |              |      |          |      |
| Methyl tert-butyl ether (MTBE) | ND        | 2.00 |                          |             |   |          |              |      |          |      |
| 1,2,4-Trimethylbenzene         | ND        | 2.00 |                          |             |   |          |              |      |          |      |
| 1,3,5-Trimethylbenzene         | ND        | 2.00 |                          |             |   |          |              |      |          |      |
| 1,2-Dichloroethane (EDC)       | ND        | 2.00 |                          |             |   |          |              |      |          |      |
| 1,2-Dibromoethane (EDB)        | ND        | 2.00 |                          |             |   |          |              |      |          |      |
| Naphthalene                    | ND        | 2.00 |                          |             |   |          |              |      |          |      |
| 1-Methylnaphthalene            | ND        | 4.00 |                          |             |   |          |              |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

R RPD outside accepted recovery limits

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

P Sample pH Not In Range

RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                             |                  |      |                |                  |      |           |  |        |              |      |
|-----------------------------|------------------|------|----------------|------------------|------|-----------|--|--------|--------------|------|
| Sample ID                   | <b>MB-25468</b>  |      | SampType:      | <b>MBLK</b>      |      | TestCode: | <b>Method 8260B/5035LOW: VOLATILES</b> |        |              |      |
| Client ID:                  | <b>PBS</b>       |      | Batch ID:      | <b>25468</b>     |      | RunNo:    | <b>34484</b>                           |        |              |      |
| Prep Date:                  | <b>5/24/2016</b> |      | Analysis Date: | <b>5/25/2016</b> |      | SeqNo:    | <b>1063541</b>                         | Units: | <b>µg/Kg</b> |      |
| Analyte                     | Result           | PQL  | SPK value      | SPK Ref Val      | %REC | LowLimit  | HighLimit                              | %RPD   | RPDLimit     | Qual |
| 2-Methylnaphthalene         | ND               | 4.00 |                |                  |      |           |  |        |              |      |
| Acetone                     | ND               | 10.0 |                |                  |      |           |  |        |              |      |
| Bromobenzene                | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Bromodichloromethane        | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Bromoform                   | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Bromomethane                | ND               | 3.00 |                |                  |      |           |  |        |              |      |
| 2-Butanone                  | ND               | 10.0 |                |                  |      |           |  |        |              |      |
| Carbon disulfide            | ND               | 10.0 |                |                  |      |           |  |        |              |      |
| Carbon tetrachloride        | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Chlorobenzene               | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Chloroethane                | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Chloroform                  | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Chloromethane               | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 2-Chlorotoluene             | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 4-Chlorotoluene             | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| cis-1,2-DCE                 | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| cis-1,3-Dichloropropene     | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 1,2-Dibromo-3-chloropropane | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Dibromochloromethane        | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Dibromomethane              | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 1,2-Dichlorobenzene         | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 1,3-Dichlorobenzene         | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 1,4-Dichlorobenzene         | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Dichlorodifluoromethane     | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 1,1-Dichloroethane          | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 1,1-Dichloroethene          | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 1,2-Dichloropropane         | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 1,3-Dichloropropane         | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 2,2-Dichloropropane         | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 1,1-Dichloropropene         | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| Hexachlorobutadiene         | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 2-Hexanone                  | ND               | 10.0 |                |                  |      |           |  |        |              |      |
| Isopropylbenzene            | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 4-Isopropyltoluene          | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| 4-Methyl-2-pentanone        | ND               | 10.0 |                |                  |      |           |  |        |              |      |
| Methylene chloride          | ND               | 3.00 |                |                  |      |           |  |        |              |      |
| n-Butylbenzene              | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| n-Propylbenzene             | ND               | 2.00 |                |                  |      |           |  |        |              |      |
| sec-Butylbenzene            | ND               | 2.00 |                |                  |      |           |  |        |              |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                             |           |      |                |             |      |           |                                 |        |          |      |
|-----------------------------|-----------|------|----------------|-------------|------|-----------|---------------------------------|--------|----------|------|
| Sample ID                   | MB-25468  |      | SampType:      | MBLK        |      | TestCode: | Method 8260B/5035LOW: VOLATILES |        |          |      |
| Client ID:                  | PBS       |      | Batch ID:      | 25468       |      | RunNo:    | 34484                           |        |          |      |
| Prep Date:                  | 5/24/2016 |      | Analysis Date: | 5/25/2016   |      | SeqNo:    | 1063541                         | Units: | µg/Kg    |      |
| Analyte                     | Result    | PQL  | SPK value      | SPK Ref Val | %REC | LowLimit  | HighLimit                       | %RPD   | RPDLimit | Qual |
| Styrene                     | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| tert-Butylbenzene           | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| 1,1,1,2-Tetrachloroethane   | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| 1,1,2,2-Tetrachloroethane   | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| Tetrachloroethene (PCE)     | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| trans-1,2-DCE               | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| trans-1,3-Dichloropropene   | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| 1,2,3-Trichlorobenzene      | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| 1,2,4-Trichlorobenzene      | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| 1,1,1-Trichloroethane       | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| 1,1,2-Trichloroethane       | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| Trichloroethene (TCE)       | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| Trichlorofluoromethane      | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| 1,2,3-Trichloropropane      | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| Vinyl chloride              | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| Xylenes, Total              | ND        | 2.00 |                |             |      |           |                                 |        |          |      |
| Surr: 1,2-Dichloroethane-d4 | 8.94      |      | 10.00          |             | 89.4 | 70        | 130                             |        |          |      |
| Surr: 4-Bromofluorobenzene  | 10.0      |      | 10.00          |             | 100  | 70        | 130                             |        |          |      |
| Surr: Dibromofluoromethane  | 8.80      |      | 10.00          |             | 88.0 | 70        | 130                             |        |          |      |
| Surr: Toluene-d8            | 9.82      |      | 10.00          |             | 98.2 | 70        | 130                             |        |          |      |

### Qualifiers:

|    |   |    |   |
|----|---|----|---|
| *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
| D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
| H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
| ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
| R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
| S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |



# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                             |           |                |           |             |                                 |          |           |      |          |      |
|-----------------------------|-----------|----------------|-----------|-------------|---------------------------------|----------|-----------|------|----------|------|
| Sample ID                   | mb-25469  | SampType:      | MBLK      | TestCode:   | EPA Method 8270C: Semivolatiles |          |           |      |          |      |
| Client ID:                  | PBS       | Batch ID:      | 25469     | RunNo:      | 34461                           |          |           |      |          |      |
| Prep Date:                  | 5/24/2016 | Analysis Date: | 5/25/2016 | SeqNo:      | 1063511                         | Units:   | mg/Kg     |      |          |      |
| Analyte                     | Result    | PQL            | SPK value | SPK Ref Val | %REC                            | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
| Acenaphthene                | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Acenaphthylene              | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Aniline                     | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Anthracene                  | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Azobenzene                  | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Benz(a)anthracene           | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Benzo(a)pyrene              | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Benzo(b)fluoranthene        | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Benzo(g,h,i)perylene        | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Benzo(k)fluoranthene        | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Benzoic acid                | ND        | 0.50           |           |             |                                 |          |           |      |          |      |
| Benzyl alcohol              | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Bis(2-chloroethoxy)methane  | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Bis(2-chloroethyl)ether     | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Bis(2-chloroisopropyl)ether | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Bis(2-ethylhexyl)phthalate  | 0.15      | 0.50           |           |             |                                 |          |           |      |          | J    |
| 4-Bromophenyl phenyl ether  | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Butyl benzyl phthalate      | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Carbazole                   | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| 4-Chloro-3-methylphenol     | ND        | 0.50           |           |             |                                 |          |           |      |          |      |
| 4-Chloroaniline             | ND        | 0.50           |           |             |                                 |          |           |      |          |      |
| 2-Chloronaphthalene         | ND        | 0.25           |           |             |                                 |          |           |      |          |      |
| 2-Chlorophenol              | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| 4-Chlorophenyl phenyl ether | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Chrysene                    | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Di-n-butyl phthalate        | 0.12      | 0.40           |           |             |                                 |          |           |      |          | J    |
| Di-n-octyl phthalate        | ND        | 0.40           |           |             |                                 |          |           |      |          |      |
| Dibenz(a,h)anthracene       | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Dibenzofuran                | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| 1,2-Dichlorobenzene         | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| 1,3-Dichlorobenzene         | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| 1,4-Dichlorobenzene         | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| 3,3'-Dichlorobenzidine      | ND        | 0.25           |           |             |                                 |          |           |      |          |      |
| Diethyl phthalate           | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| Dimethyl phthalate          | ND        | 0.20           |           |             |                                 |          |           |      |          |      |
| 2,4-Dichlorophenol          | ND        | 0.40           |           |             |                                 |          |           |      |          |      |
| 2,4-Dimethylphenol          | ND        | 0.30           |           |             |                                 |          |           |      |          |      |
| 4,6-Dinitro-2-methylphenol  | 0.087     | 0.40           |           |             |                                 |          |           |      |          | J    |
| 2,4-Dinitrophenol           | ND        | 0.50           |           |             |                                 |          |           |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                            |           |      |                |             |      |           |                                 |      |              |      |
|----------------------------|-----------|------|----------------|-------------|------|-----------|---------------------------------|------|--------------|------|
| Sample ID                  | mb-25469  |      | SampType:      | MBLK        |      | TestCode: | EPA Method 8270C: Semivolatiles |      |              |      |
| Client ID:                 | PBS       |      | Batch ID:      | 25469       |      | RunNo:    | 34461                           |      |              |      |
| Prep Date:                 | 5/24/2016 |      | Analysis Date: | 5/25/2016   |      | SeqNo:    | 1063511                         |      | Units: mg/Kg |      |
| Analyte                    | Result    | PQL  | SPK value      | SPK Ref Val | %REC | LowLimit  | HighLimit                       | %RPD | RPDLimit     | Qual |
| 2,4-Dinitrotoluene         | ND        | 0.50 |                |             |      |           |                                 |      |              |      |
| 2,6-Dinitrotoluene         | ND        | 0.50 |                |             |      |           |                                 |      |              |      |
| Fluoranthene               | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Fluorene                   | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Hexachlorobenzene          | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Hexachlorobutadiene        | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Hexachlorocyclopentadiene  | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Hexachloroethane           | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Indeno(1,2,3-cd)pyrene     | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| 1-Methylnaphthalene        | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| 2-Methylnaphthalene        | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| 2-Methylphenol             | ND        | 0.40 |                |             |      |           |                                 |      |              |      |
| 3+4-Methylphenol           | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| N-Nitrosodi-n-propylamine  | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| N-Nitrosodiphenylamine     | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Naphthalene                | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| 2-Nitroaniline             | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| 3-Nitroaniline             | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| 4-Nitroaniline             | ND        | 0.40 |                |             |      |           |                                 |      |              |      |
| Nitrobenzene               | ND        | 0.40 |                |             |      |           |                                 |      |              |      |
| 2-Nitrophenol              | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| 4-Nitrophenol              | ND        | 0.25 |                |             |      |           |                                 |      |              |      |
| Pentachlorophenol          | ND        | 0.40 |                |             |      |           |                                 |      |              |      |
| Phenanthrene               | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Phenol                     | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Pyrene                     | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Pyridine                   | ND        | 0.40 |                |             |      |           |                                 |      |              |      |
| 1,2,4-Trichlorobenzene     | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| 2,4,5-Trichlorophenol      | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| 2,4,6-Trichlorophenol      | ND        | 0.20 |                |             |      |           |                                 |      |              |      |
| Surr: 2-Fluorophenol       | 2.0       |      | 3.330          |             | 60.1 | 28.3      | 102                             |      |              |      |
| Surr: Phenol-d5            | 2.2       |      | 3.330          |             | 65.9 | 35.7      | 103                             |      |              |      |
| Surr: 2,4,6-Tribromophenol | 2.5       |      | 3.330          |             | 75.8 | 35.2      | 108                             |      |              |      |
| Surr: Nitrobenzene-d5      | 1.1       |      | 1.670          |             | 68.4 | 24        | 118                             |      |              |      |
| Surr: 2-Fluorobiphenyl     | 1.2       |      | 1.670          |             | 73.2 | 35.4      | 111                             |      |              |      |
| Surr: 4-Terphenyl-d14      | 1.1       |      | 1.670          |             | 64.4 | 15        | 91.7                            |      |              |      |

### Qualifiers:

|    |   |    |   |
|----|---|----|---|
| *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
| D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
| H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
| ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
| R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
| S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                            |           |      |                          |             |      |   |           |              |          |      |
|----------------------------|-----------|------|--------------------------|-------------|------|---|-----------|--------------|----------|------|
| Sample ID                  | Ics-25469 |      | SampType: LCS            |             |      | TestCode: EPA Method 8270C: Semivolatiles |           |              |          |      |
| Client ID:                 | LCSS      |      | Batch ID: 25469          |             |      | RunNo: 34461                              |           |              |          |      |
| Prep Date:                 | 5/24/2016 |      | Analysis Date: 5/25/2016 |             |      | SeqNo: 1063512                            |           | Units: mg/Kg |          |      |
| Analyte                    | Result    | PQL  | SPK value                | SPK Ref Val | %REC | LowLimit                                  | HighLimit | %RPD         | RPDLimit | Qual |
| Acenaphthene               | 1.2       | 0.20 | 1.670                    | 0           | 70.0 | 45.8                                      | 99.8      |              |          |      |
| 4-Chloro-3-methylphenol    | 2.2       | 0.50 | 3.330                    | 0           | 66.5 | 51.5                                      | 103       |              |          |      |
| 2-Chlorophenol             | 1.8       | 0.20 | 3.330                    | 0           | 54.1 | 46.5                                      | 105       |              |          |      |
| 1,4-Dichlorobenzene        | 0.86      | 0.20 | 1.670                    | 0           | 51.3 | 45.5                                      | 103       |              |          |      |
| 2,4-Dinitrotoluene         | 1.1       | 0.50 | 1.670                    | 0           | 66.9 | 36  | 87.2      |              |          |      |
| N-Nitrosodi-n-propylamine  | 0.80      | 0.20 | 1.670                    | 0           | 48.1 | 47.3                                      | 104       |              |          |      |
| 4-Nitrophenol              | 2.2       | 0.25 | 3.330                    | 0           | 65.7 | 47.3                                      | 95.3      |              |          |      |
| Pentachlorophenol          | 1.9       | 0.40 | 3.330                    | 0           | 56.6 | 38.7                                      | 89.3      |              |          |      |
| Phenol                     | 1.8       | 0.20 | 3.330                    | 0           | 54.5 | 47.8                                      | 106       |              |          |      |
| Pyrene                     | 1.2       | 0.20 | 1.670                    | 0           | 72.6 | 33.4                                      | 105       |              |          |      |
| 1,2,4-Trichlorobenzene     | 1.0       | 0.20 | 1.670                    | 0           | 61.4 | 50.4                                      | 115       |              |          |      |
| Surr: 2-Fluorophenol       | 1.6       |      | 3.330                    |             | 47.4 | 28.3                                      | 102       |              |          |      |
| Surr: Phenol-d5            | 1.9       |      | 3.330                    |             | 56.0 | 35.7                                      | 103       |              |          |      |
| Surr: 2,4,6-Tribromophenol | 2.2       |      | 3.330                    |             | 66.3 | 35.2                                      | 108       |              |          |      |
| Surr: Nitrobenzene-d5      | 1.0       |      | 1.670                    |             | 59.7 | 24  | 118       |              |          |      |
| Surr: 2-Fluorobiphenyl     | 1.2       |      | 1.670                    |             | 70.2 | 35.4                                      | 111       |              |          |      |
| Surr: 4-Terphenyl-d14      | 0.74      |      | 1.670                    |             | 44.5 | 15  | 91.7      |              |          |      |

### Qualifiers:

|    |   |    |   |
|----|---|----|---|
| *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
| D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
| H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
| ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
| R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
| S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|            |           |       |                |             |      |           |                          |      |          |       |  |
|------------|-----------|-------|----------------|-------------|------|-----------|--------------------------|------|----------|-------|--|
| Sample ID  | MB-25509  |       | SampType:      | MBLK        |      | TestCode: | EPA Method 7471: Mercury |      |          |       |  |
| Client ID: | PBS       |       | Batch ID:      | 25509       |      | RunNo:    | 34523                    |      |          |       |  |
| Prep Date: | 5/25/2016 |       | Analysis Date: | 5/26/2016   |      | SeqNo:    | 1064862                  |      | Units:   | mg/Kg |  |
| Analyte    | Result    | PQL   | SPK value      | SPK Ref Val | %REC | LowLimit  | HighLimit                | %RPD | RPDLimit | Qual  |  |
| Mercurv    | ND        | 0.033 |                |             |      |           |                          |      |          |       |  |

|            |           |       |                          |             |                                    |          |              |      |          |      |
|------------|-----------|-------|--------------------------|-------------|------------------------------------|----------|--------------|------|----------|------|
| Sample ID  | LCS-25509 |       | SampType: LCS            |             | TestCode: EPA Method 7471: Mercury |          |              |      |          |      |
| Client ID: | LCSS      |       | Batch ID: 25509          |             | RunNo: 34523                       |          |              |      |          |      |
| Prep Date: | 5/25/2016 |       | Analysis Date: 5/26/2016 |             | SeqNo: 1064863                     |          | Units: mg/Kg |      |          |      |
| Analyte    | Result    | PQL   | SPK value                | SPK Ref Val | %REC                               | LowLimit | HighLimit    | %RPD | RPDLimit | Qual |
| Mercury    | 0.16      | 0.033 | 0.1667                   | 0           | 98.8                               | 80       | 120          |      |          |      |

|            |                     |       |                          |             |                                    |          |                  |      |          |      |
|------------|---------------------|-------|--------------------------|-------------|------------------------------------|----------|------------------|------|----------|------|
| Sample ID  | 1605943-001BMS      |       | SampType: MS             |             | TestCode: EPA Method 7471: Mercury |          |                  |      |          |      |
| Client ID: | SWMU 10-21 (2-2.5') |       | Batch ID: 25509          |             | RunNo: 34523                       |          |                  |      |          |      |
| Prep Date: | 5/25/2016           |       | Analysis Date: 5/26/2016 |             | SeqNo: 1064865                     |          | Units: mg/Kg-dry |      |          |      |
| Analyte    | Result              | PQL   | SPK value                | SPK Ref Val | %REC                               | LowLimit | HighLimit        | %RPD | RPDLimit | Qual |
| Mercury    | 0.23                | 0.036 | 0.1803                   | 0.05558     | 98.8                               | 75       | 125              |      |          |      |

|            |                     |       |                          |             |                                    |          |                  |      |          |      |
|------------|---------------------|-------|--------------------------|-------------|------------------------------------|----------|------------------|------|----------|------|
| Sample ID  | 1605943-001BMSD     |       | SampType: MSD            |             | TestCode: EPA Method 7471: Mercury |          |                  |      |          |      |
| Client ID: | SWMU 10-21 (2-2.5') |       | Batch ID: 25509          |             | RunNo: 34523                       |          |                  |      |          |      |
| Prep Date: | 5/25/2016           |       | Analysis Date: 5/26/2016 |             | SeqNo: 1064866                     |          | Units: mg/Kg-dry |      |          |      |
| Analyte    | Result              | PQL   | SPK value                | SPK Ref Val | %REC                               | LowLimit | HighLimit        | %RPD | RPDLimit | Qual |
| Mercury    | 0.23                | 0.036 | 0.1817                   | 0.05558     | 93.9                               | 75       | 125              | 3.25 | 20       |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|            |                  |      |                |                  |      |           |                                      |      |                     |      |
|------------|------------------|------|----------------|------------------|------|-----------|--------------------------------------|------|---------------------|------|
| Sample ID  | <b>MB-25508</b>  |      | SampType:      | <b>MBLK</b>      |      | TestCode: | <b>EPA Method 6010B: Soil Metals</b> |      |                     |      |
| Client ID: | <b>PBS</b>       |      | Batch ID:      | <b>25508</b>     |      | RunNo:    | <b>34511</b>                         |      |                     |      |
| Prep Date: | <b>5/25/2016</b> |      | Analysis Date: | <b>5/26/2016</b> |      | SeqNo:    | <b>1064307</b>                       |      | Units: <b>mg/Kg</b> |      |
| Analyte    | Result           | PQL  | SPK value      | SPK Ref Val      | %REC | LowLimit  | HighLimit                            | %RPD | RPDLimit            | Qual |
| Antimony   | ND               | 2.5  |                |                  |      |           |                                      |      |                     |      |
| Arsenic    | ND               | 2.5  |                |                  |      |           |                                      |      |                     |      |
| Barium     | ND               | 0.10 |                |                  |      |           |                                      |      |                     |      |
| Beryllium  | ND               | 0.15 |                |                  |      |           |                                      |      |                     |      |
| Cadmium    | ND               | 0.10 |                |                  |      |           |                                      |      |                     |      |
| Chromium   | ND               | 0.30 |                |                  |      |           |                                      |      |                     |      |
| Cobalt     | ND               | 0.30 |                |                  |      |           |                                      |      |                     |      |
| Iron       | 1.7              | 2.5  |                |                  |      |           |                                      |      |                     | J    |
| Lead       | ND               | 0.25 |                |                  |      |           |                                      |      |                     |      |
| Manganese  | ND               | 0.10 |                |                  |      |           |                                      |      |                     |      |
| Nickel     | ND               | 0.50 |                |                  |      |           |                                      |      |                     |      |
| Selenium   | ND               | 2.5  |                |                  |      |           |                                      |      |                     |      |
| Silver     | ND               | 0.25 |                |                  |      |           |                                      |      |                     |      |
| Vanadium   | ND               | 2.5  |                |                  |      |           |                                      |      |                     |      |
| Zinc       | ND               | 2.5  |                |                  |      |           |                                      |      |                     |      |

|            |                  |      |                |                  |      |           |                                      |      |                     |      |
|------------|------------------|------|----------------|------------------|------|-----------|--------------------------------------|------|---------------------|------|
| Sample ID  | <b>LCS-25508</b> |      | SampType:      | <b>LCS</b>       |      | TestCode: | <b>EPA Method 6010B: Soil Metals</b> |      |                     |      |
| Client ID: | <b>LCSS</b>      |      | Batch ID:      | <b>25508</b>     |      | RunNo:    | <b>34511</b>                         |      |                     |      |
| Prep Date: | <b>5/25/2016</b> |      | Analysis Date: | <b>5/26/2016</b> |      | SeqNo:    | <b>1064308</b>                       |      | Units: <b>mg/Kg</b> |      |
| Analyte    | Result           | PQL  | SPK value      | SPK Ref Val      | %REC | LowLimit  | HighLimit                            | %RPD | RPDLimit            | Qual |
| Antimony   | 25               | 2.5  | 25.00          | 0                | 99.8 | 80        | 120                                  |      |                     |      |
| Arsenic    | 26               | 2.5  | 25.00          | 0                | 104  | 80        | 120                                  |      |                     |      |
| Barium     | 25               | 0.10 | 25.00          | 0                | 99.1 | 80        | 120                                  |      |                     |      |
| Beryllium  | 26               | 0.15 | 25.00          | 0                | 103  | 80        | 120                                  |      |                     |      |
| Cadmium    | 25               | 0.10 | 25.00          | 0                | 101  | 80        | 120                                  |      |                     |      |
| Chromium   | 25               | 0.30 | 25.00          | 0                | 100  | 80        | 120                                  |      |                     |      |
| Cobalt     | 24               | 0.30 | 25.00          | 0                | 95.6 | 80        | 120                                  |      |                     |      |
| Iron       | 26               | 2.5  | 25.00          | 0                | 105  | 80        | 120                                  |      |                     |      |
| Lead       | 25               | 0.25 | 25.00          | 0                | 98.3 | 80        | 120                                  |      |                     |      |
| Manganese  | 25               | 0.10 | 25.00          | 0                | 99.6 | 80        | 120                                  |      |                     |      |
| Nickel     | 25               | 0.50 | 25.00          | 0                | 99.1 | 80        | 120                                  |      |                     |      |
| Selenium   | 25               | 2.5  | 25.00          | 0                | 99.6 | 80        | 120                                  |      |                     |      |
| Silver     | 5.0              | 0.25 | 5.000          | 0                | 101  | 80        | 120                                  |      |                     |      |
| Vanadium   | 26               | 2.5  | 25.00          | 0                | 103  | 80        | 120                                  |      |                     |      |
| Zinc       | 25               | 2.5  | 25.00          | 0                | 99.1 | 80        | 120                                  |      |                     |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

**Client:** Western Refining Company

**Project:** SWMU 10

|            |                     |      |                          |             |   |          |                  |      |          |      |
|------------|---------------------|------|--------------------------|-------------|---|----------|------------------|------|----------|------|
| Sample ID  | 1605943-001BMS      |      | SampType: MS             |             | TestCode: EPA Method 6010B: Soil Metals |          |                  |      |          |      |
| Client ID: | SWMU 10-21 (2-2.5') |      | Batch ID: 25508          |             | RunNo: 34511                            |          |                  |      |          |      |
| Prep Date: | 5/25/2016           |      | Analysis Date: 5/26/2016 |             | SeqNo: 1064368                          |          | Units: mg/Kg-dry |      |          |      |
| Analyte    | Result              | PQL  | SPK value                | SPK Ref Val | %REC                                    | LowLimit | HighLimit        | %RPD | RPDLimit | Qual |
| Arsenic    | 28                  | 2.7  | 27.29                    | 3.337       | 90.8                                    | 75       | 125              |      |          |      |
| Beryllium  | 25                  | 0.16 | 27.29                    | 0.7180      | 89.7                                    | 75       | 125              |      |          |      |
| Cadmium    | 24                  | 0.11 | 27.29                    | 0           | 88.8                                    | 75       | 125              |      |          |      |
| Chromium   | 35                  | 0.33 | 27.29                    | 9.159       | 95.8                                    | 75       | 125              |      |          |      |
| Cobalt     | 25                  | 0.33 | 27.29                    | 3.682       | 76.6                                    | 75       | 125              |      |          |      |
| Silver     | 4.8                 | 0.27 | 5.459                    | 0           | 88.0                                    | 75       | 125              |      |          |      |
| Vanadium   | 47                  | 2.7  | 27.29                    | 16.05       | 112                                     | 75       | 125              |      |          |      |
| Zinc       | 50                  | 2.7  | 27.29                    | 16.14       | 123                                     | 75       | 125              |      |          |      |

|            |                     |      |                          |             |   |          |                  |       |          |      |
|------------|---------------------|------|--------------------------|-------------|---|----------|------------------|-------|----------|------|
| Sample ID  | 1605943-001BMSD     |      | SampType: MSD            |             | TestCode: EPA Method 6010B: Soil Metals |          |                  |       |          |      |
| Client ID: | SWMU 10-21 (2-2.5') |      | Batch ID: 25508          |             | RunNo: 34511                            |          |                  |       |          |      |
| Prep Date: | 5/25/2016           |      | Analysis Date: 5/26/2016 |             | SeqNo: 1064369                          |          | Units: mg/Kg-dry |       |          |      |
| Analyte    | Result              | PQL  | SPK value                | SPK Ref Val | %REC                                    | LowLimit | HighLimit        | %RPD  | RPDLimit | Qual |
| Arsenic    | 29                  | 2.7  | 26.83                    | 3.337       | 96.3                                    | 75       | 125              | 3.65  | 20       |      |
| Beryllium  | 24                  | 0.16 | 26.83                    | 0.7180      | 88.6                                    | 75       | 125              | 2.80  | 20       |      |
| Cadmium    | 24                  | 0.11 | 26.83                    | 0           | 91.2                                    | 75       | 125              | 1.00  | 20       |      |
| Chromium   | 30                  | 0.32 | 26.83                    | 9.159       | 76.9                                    | 75       | 125              | 16.9  | 20       |      |
| Cobalt     | 23                  | 0.32 | 26.83                    | 3.682       | 72.4                                    | 75       | 125              | 6.20  | 20       | S    |
| Silver     | 4.8                 | 0.27 | 5.367                    | 0           | 89.2                                    | 75       | 125              | 0.310 | 20       |      |
| Vanadium   | 40                  | 2.7  | 26.83                    | 16.05       | 88.1                                    | 75       | 125              | 15.8  | 20       |      |
| Zinc       | 35                  | 2.7  | 26.83                    | 16.14       | 70.7                                    | 75       | 125              | 34.4  | 20       | RS   |

|            |                     |      |                          |             |   |          |                  |      |          |      |
|------------|---------------------|------|--------------------------|-------------|---|----------|------------------|------|----------|------|
| Sample ID  | 1605943-001BMS      |      | SampType: MS             |             | TestCode: EPA Method 6010B: Soil Metals |          |                  |      |          |      |
| Client ID: | SWMU 10-21 (2-2.5') |      | Batch ID: 25508          |             | RunNo: 34558                            |          |                  |      |          |      |
| Prep Date: | 5/25/2016           |      | Analysis Date: 5/31/2016 |             | SeqNo: 1065752                          |          | Units: mg/Kg-dry |      |          |      |
| Analyte    | Result              | PQL  | SPK value                | SPK Ref Val | %REC                                    | LowLimit | HighLimit        | %RPD | RPDLimit | Qual |
| Antimony   | 8.7                 | 2.7  | 27.29                    | 0           | 31.9                                    | 75       | 125              |      |          | S    |
| Lead       | 23                  | 0.27 | 27.29                    | 2.127       | 74.7                                    | 75       | 125              |      |          | S    |
| Nickel     | 27                  | 0.55 | 27.29                    | 6.041       | 78.0                                    | 75       | 125              |      |          |      |
| Selenium   | 21                  | 2.7  | 27.29                    | 0           | 75.9                                    | 75       | 125              |      |          |      |

|            |                     |      |                |             |      |           |                               |      |                  |      |  |
|------------|---------------------|------|----------------|-------------|------|-----------|-------------------------------|------|------------------|------|--|
| Sample ID  | 1605943-001BMSD     |      | SampType:      | MSD         |      | TestCode: | EPA Method 6010B: Soil Metals |      |                  |      |  |
| Client ID: | SWMU 10-21 (2-2.5') |      | Batch ID:      | 25508       |      | RunNo:    | 34558                         |      |                  |      |  |
| Prep Date: | 5/25/2016           |      | Analysis Date: | 5/31/2016   |      | SeqNo:    | 1065753                       |      | Units: mg/Kg-dry |      |  |
| Analyte    | Result              | PQL  | SPK value      | SPK Ref Val | %REC | LowLimit  | HighLimit                     | %RPD | RPDLimit         | Qual |  |
| Antimony   | 4.5                 | 2.7  | 26.83          | 0           | 16.9 | 75        | 125                           | 63.2 | 20               | RS   |  |
| Lead       | 21                  | 0.27 | 26.83          | 2.127       | 71.1 | 75        | 125                           | 5.99 | 20               | S    |  |

### Qualifiers:

- |   |   |
|---|---|
| * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
| D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
| H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
| ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
| R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
| S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

Client: Western Refining Company

Project: SWMU 10

|            |                     |                |           |             |                               |          |           |      |          |      |
|------------|---------------------|----------------|-----------|-------------|-------------------------------|----------|-----------|------|----------|------|
| Sample ID  | 1605943-001BMSD     | SampType:      | MSD       | TestCode:   | EPA Method 6010B: Soil Metals |          |           |      |          |      |
| Client ID: | SWMU 10-21 (2-2.5') | Batch ID:      | 25508     | RunNo:      | 34558                         |          |           |      |          |      |
| Prep Date: | 5/25/2016           | Analysis Date: | 5/31/2016 | SeqNo:      | 1065753                       | Units:   | mg/Kg-dry |      |          |      |
| Analyte    | Result              | PQL            | SPK value | SPK Ref Val | %REC                          | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
| Nickel     | 25                  | 0.54           | 26.83     | 6.041       | 71.1                          | 75       | 125       | 8.39 | 20       | S    |
| Selenium   | 19                  | 2.7            | 26.83     | 0           | 69.3                          | 75       | 125       | 10.9 | 20       | S    |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605943

28-Jun-16

**Client:** Western Refining Company

**Project:** SWMU 10

| Sample ID <b>rb</b>           | SampType: <b>MBLK</b>           |     |           |             | TestCode: <b>EPA Method 8015D Mod: Gasoline Range</b> |          |                     |      |          |      |
|-------------------------------|---------------------------------|-----|-----------|-------------|---|----------|---------------------|------|----------|------|
| Client ID: <b>PBS</b>         | Batch ID: <b>A34454</b>         |     |           |             | RunNo: <b>34454</b>                                   |          |                     |      |          |      |
| Prep Date:                    | Analysis Date: <b>5/24/2016</b> |     |           |             | SeqNo: <b>1062581</b>                                 |          | Units: <b>mg/Kg</b> |      |          |      |
| Analyte                       | Result                          | PQL | SPK value | SPK Ref Val | %REC  | LowLimit | HighLimit           | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO) | ND                              | 5.0 |           |             |   |          |                     |      |          |      |
| Surr: BFB                     | 510                             |     | 500.0     |             | 101   | 70       | 130                 |      |          |      |

| Sample ID <b>2.5ug gro lcs</b> | SampType: <b>LCS</b>            |     |           |             | TestCode: <b>EPA Method 8015D Mod: Gasoline Range</b> |          |                     |      |          |      |
|--------------------------------|---------------------------------|-----|-----------|-------------|---|----------|---------------------|------|----------|------|
| Client ID: <b>LCSS</b>         | Batch ID: <b>A34454</b>         |     |           |             | RunNo: <b>34454</b>                                   |          |                     |      |          |      |
| Prep Date:                     | Analysis Date: <b>5/24/2016</b> |     |           |             | SeqNo: <b>1062582</b>                                 |          | Units: <b>mg/Kg</b> |      |          |      |
| Analyte                        | Result                          | PQL | SPK value | SPK Ref Val | %REC  | LowLimit | HighLimit           | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO)  | 26                              | 5.0 | 25.00     | 0           | 103   | 62.9     | 123                 |      |          |      |
| Surr: BFB                      | 500                             |     | 500.0     |             | 99.4  | 70       | 130                 |      |          |      |

| Sample ID <b>1605943-001bms</b>       | SampType: <b>MS</b>             |     |           |             | TestCode: <b>EPA Method 8015D Mod: Gasoline Range</b> |          |                         |      |          |      |
|---------------------------------------|---------------------------------|-----|-----------|-------------|---|----------|-------------------------|------|----------|------|
| Client ID: <b>SWMU 10-21 (2-2.5')</b> | Batch ID: <b>A34454</b>         |     |           |             | RunNo: <b>34454</b>                                   |          |                         |      |          |      |
| Prep Date:                            | Analysis Date: <b>5/24/2016</b> |     |           |             | SeqNo: <b>1062584</b>                                 |          | Units: <b>mg/Kg-dry</b> |      |          |      |
| Analyte                               | Result                          | PQL | SPK value | SPK Ref Val | %REC  | LowLimit | HighLimit               | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO)         | 24                              | 4.1 | 20.41     | 1.968       | 106   | 52.3     | 132                     |      |          |      |
| Surr: BFB                             | 410                             |     | 408.2     |             | 101   | 70       | 130                     |      |          |      |

| Sample ID <b>1605943-001bmsd</b>      | SampType: <b>MSD</b>            |     |           |             | TestCode: <b>EPA Method 8015D Mod: Gasoline Range</b> |          |                         |      |          |      |
|---------------------------------------|---------------------------------|-----|-----------|-------------|---|----------|-------------------------|------|----------|------|
| Client ID: <b>SWMU 10-21 (2-2.5')</b> | Batch ID: <b>A34454</b>         |     |           |             | RunNo: <b>34454</b>                                   |          |                         |      |          |      |
| Prep Date:                            | Analysis Date: <b>5/24/2016</b> |     |           |             | SeqNo: <b>1062585</b>                                 |          | Units: <b>mg/Kg-dry</b> |      |          |      |
| Analyte                               | Result                          | PQL | SPK value | SPK Ref Val | %REC  | LowLimit | HighLimit               | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO)         | 22                              | 4.1 | 20.41     | 1.968       | 99.4  | 52.3     | 132                     | 6.01 | 20       |      |
| Surr: BFB                             | 390                             |     | 408.2     |             | 96.6  | 70       | 130                     | 0    | 0        |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified





Hall Environmental Analysis Laboratory  
4901 Hawkins NE  
Albuquerque, NM 87109  
TEL: 505-345-3975 FAX: 505-345-4107  
Website: www.hallenvironmental.com

## Sample Log-In Check List

Client Name: Western Refining Gallup

Work Order Number: 1605943

RcptNo: 1

Received by/date:

AG 05/19/16

Logged By: Lindsay Mangin

5/19/2016 4:30:00 PM

*Judy Mangin*

Completed By: Lindsay Mangin

5/20/2016 8:48:48 AM

*Judy Mangin*

Reviewed By:

*JM* 05/23/16

### Chain of Custody

1. Custody seals intact on sample bottles? Yes ☐ No ☐ Not Present ☒
2. Is Chain of Custody complete? Yes ☒ No ☐ Not Present ☐
3. How was the sample delivered? Courier

### Log In

4. Was an attempt made to cool the samples? Yes ☒ No ☐ NA ☐
5. Were all samples received at a temperature of  $>0^{\circ}\text{C}$  to  $6.0^{\circ}\text{C}$ ? Yes ☒ No ☐ NA ☐
6. Sample(s) in proper container(s)? Yes ☒ No ☐
7. Sufficient sample volume for indicated test(s)? Yes ☒ No ☐
8. Are samples (except VOA and ONG) properly preserved? Yes ☒ No ☐
9. Was preservative added to bottles? Yes ☐ No ☒ NA ☐
10. VOA vials have zero headspace? Yes ☐ No ☐ No VOA Vials ☒
11. Were any sample containers received broken? Yes ☐ No ☒
12. Does paperwork match bottle labels?  
(Note discrepancies on chain of custody) Yes ☒ No ☐
13. Are matrices correctly identified on Chain of Custody? Yes ☒ No ☐
14. Is it clear what analyses were requested? Yes ☒ No ☐
15. Were all holding times able to be met?  
(If no, notify customer for authorization.) Yes ☒ No ☐

# of preserved  
bottles checked  
for pH:

(<2 or >12 unless noted)

Adjusted? \_\_\_\_\_

Checked by: \_\_\_\_\_

### Special Handling (if applicable)

16. Was client notified of all discrepancies with this order? Yes ☐ No ☐ NA ☒

Person Notified: \_\_\_\_\_

Date: \_\_\_\_\_

By Whom: \_\_\_\_\_

Via: ☐ eMail ☐ Phone ☐ Fax ☐ In Person

Regarding: \_\_\_\_\_

Client Instructions: \_\_\_\_\_

17. Additional remarks:

### 18. Cooler Information

| Cooler No | Temp $^{\circ}\text{C}$ | Condition | Seal Intact | Seal No | Seal Date | Signed By |
|-----------|-------------------------|-----------|-------------|---------|-----------|-----------|
| 1         | 1.0                     | Good      | Yes         |         |           |           |





WESTERN REFINING SOUTHWEST, INC.  
GALLUP REFINERY - SWMU 10  
METALS AND CYANIDE ANALYSES FOR SOIL SAMPLES

| Analyte     | Analytical Method             |
|-------------|-------------------------------|
| Antimony    | SW-846 method 6010/6020       |
| Arsenic     | SW-846 method 6010/6020       |
| Barium      | SW-846 method 6010/6020       |
| Beryllium   | SW-846 method 6010/6020       |
| Cadmium     | SW-846 method 6010/6020       |
| Chromium    | SW-846 method 6010/6020       |
| Chromium VI | SW-846 Method 3060A           |
| Cobalt      | SW-846 method 6010/6020       |
| Cyanide     | SW-846 method 335.4/335.2 mod |
| Lead        | SW-846 method 6010/6020       |
| Mercury     | SW-846 method 7470/7471       |
| Nickel      | SW-846 method 6010/6020       |
| Selenium    | SW-846 method 6010/6020       |
| Silver      | SW-846 method 6010/6020       |
| Vanadium    | SW-846 method 6010/6020       |
| Zinc        | SW-846 method 6010/6020       |
| Iron        | SW-846 method 6010/6020       |
| Manganese   | SW-846 method 6010/6020       |



Hall Environmental Analysis Laboratory  
4901 Hawkins NE  
Albuquerque, NM 87109  
TEL: 505-345-3975 FAX: 505-345-4107  
Website: [www.hallenvironmental.com](http://www.hallenvironmental.com)

June 30, 2016

Ed Riege

Western Refining Company

Rt. 3 Box 7

Gallup, NM 87301

TEL: (505) 722-0231

FAX

RE: SWMU 10

OrderNo.: 1605998

Dear Ed Riege:

Hall Environmental Analysis Laboratory received 3 sample(s) on 5/20/2016 for the analyses presented in the following report.

These were analyzed according to EPA procedures or equivalent. To access our accredited tests please go to [www.hallenvironmental.com](http://www.hallenvironmental.com) or the state specific web sites. In order to properly interpret your results it is imperative that you review this report in its entirety. See the sample checklist and/or the Chain of Custody for information regarding the sample receipt temperature and preservation. Data qualifiers or a narrative will be provided if the sample analysis or analytical quality control parameters require a flag. When necessary, data qualifiers are provided on both the sample analysis report and the QC summary report, both sections should be reviewed. All samples are reported, as received, unless otherwise indicated. Lab measurement of analytes considered field parameters that require analysis within 15 minutes of sampling such as pH and residual chlorine are qualified as being analyzed outside of the recommended holding time.

Please don't hesitate to contact HEAL for any additional information or clarifications.

ADHS Cert #AZ0682 -- NMED-DWB Cert #NM9425 -- NMED-Micro Cert #NM0190

Sincerely,

A handwritten signature in black ink, appearing to read 'Andy Freeman', is written over a horizontal line.

Andy Freeman

Laboratory Manager

4901 Hawkins NE

Albuquerque, NM 87109

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-20-GW

Project: SWMU 10

Collection Date: 5/20/2016 10:45:00 AM

Lab ID: 1605998-001

Matrix: AQUEOUS

Received Date: 5/20/2016 3:00:00 PM

| Analyses                                  | Result | MDL     | PQL    | Qual | Units | DF  | Date Analyzed         | Batch ID |
|---|--------|---------|--------|------|-------|-----|-----------------------|----------|
| <b>EPA METHOD 8015M/D: DIESEL RANGE</b>   |        |         |        |      |       |     | Analyst: TOM          |          |
| Diesel Range Organics (DRO)               | 440    | 6.9     | 10     |      | mg/L  | 10  | 5/26/2016 3:40:44 PM  | 25494    |
| Motor Oil Range Organics (MRO)            | ND     | 50      | 50     |      | mg/L  | 10  | 5/26/2016 3:40:44 PM  | 25494    |
| Surr: DNOP                                | 0      | 0       | 70-141 | S    | %Rec  | 10  | 5/26/2016 3:40:44 PM  | 25494    |
| <b>EPA METHOD 300.0: ANIONS</b>           |        |         |        |      |       |     | Analyst: LGT          |          |
| Fluoride                                  | ND     | 0.11    | 0.50   |      | mg/L  | 5   | 5/23/2016 5:26:04 PM  | R34434   |
| Chloride                                  | 1600   | 2.7     | 50     |      | mg/L  | 100 | 5/25/2016 3:02:36 PM  | R34487   |
| Sulfate                                   | 260    | 1.3     | 10     |      | mg/L  | 20  | 5/23/2016 5:38:29 PM  | R34434   |
| <b>EPA METHOD 200.7: DISSOLVED METALS</b> |        |         |        |      |       |     | Analyst: ELS          |          |
| Barium                                    | 0.11   | 0.0013  | 0.0020 |      | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| Beryllium                                 | ND     | 0.00031 | 0.0020 |      | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| Cadmium                                   | ND     | 0.00075 | 0.0020 |      | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| Chromium                                  | ND     | 0.0018  | 0.0060 |      | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| Cobalt                                    | 0.0074 | 0.00074 | 0.0060 |      | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| Iron                                      | 0.42   | 0.0091  | 0.020  | *    | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| Manganese                                 | 1.2    | 0.0016  | 0.010  | *    | mg/L  | 5   | 5/25/2016 1:49:50 PM  | C34463   |
| Nickel                                    | 0.019  | 0.0024  | 0.010  |      | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| Silver                                    | ND     | 0.0028  | 0.0050 |      | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| Vanadium                                  | 0.010  | 0.0013  | 0.050  | J    | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| Zinc                                      | 0.046  | 0.0028  | 0.010  |      | mg/L  | 1   | 5/25/2016 1:47:55 PM  | C34463   |
| <b>EPA METHOD 200.7: METALS</b>           |        |         |        |      |       |     | Analyst: ELS          |          |
| Barium                                    | 0.26   | 0.0013  | 0.0020 |      | mg/L  | 1   | 5/25/2016 11:04:43 AM | 25486    |
| Beryllium                                 | 0.0011 | 0.00031 | 0.0020 | J    | mg/L  | 1   | 5/25/2016 11:04:43 AM | 25486    |
| Cadmium                                   | ND     | 0.00075 | 0.0020 |      | mg/L  | 1   | 5/25/2016 11:04:43 AM | 25486    |
| Chromium                                  | 0.030  | 0.0022  | 0.0060 |      | mg/L  | 1   | 5/25/2016 11:04:43 AM | 25486    |
| Cobalt                                    | 0.011  | 0.00085 | 0.0060 |      | mg/L  | 1   | 5/25/2016 11:04:43 AM | 25486    |
| Iron                                      | 7.1    | 0.091   | 0.20   | *    | mg/L  | 10  | 6/1/2016 6:14:15 PM   | 25486    |
| Manganese                                 | 1.4    | 0.015   | 0.020  | *    | mg/L  | 10  | 5/27/2016 12:19:04 PM | 25486    |
| Nickel                                    | 0.026  | 0.0024  | 0.010  |      | mg/L  | 1   | 5/25/2016 11:04:43 AM | 25486    |
| Silver                                    | ND     | 0.0028  | 0.0050 |      | mg/L  | 1   | 5/25/2016 11:04:43 AM | 25486    |
| Vanadium                                  | 0.024  | 0.0015  | 0.050  | J    | mg/L  | 1   | 5/25/2016 11:04:43 AM | 25486    |
| Zinc                                      | 0.043  | 0.0039  | 0.010  |      | mg/L  | 1   | 5/25/2016 11:04:43 AM | 25486    |
| <b>EPA 200.8: DISSOLVED METALS</b>        |        |         |        |      |       |     | Analyst: DBD          |          |
| Antimony                                  | ND     | 0.0024  | 0.0050 |      | mg/L  | 5   | 5/23/2016 5:07:26 PM  | C34426   |
| Arsenic                                   | ND     | 0.050   | 0.050  |      | mg/L  | 50  | 5/23/2016 6:56:47 PM  | C34426   |
| Lead                                      | ND     | 0.00084 | 0.0025 |      | mg/L  | 5   | 5/23/2016 5:07:26 PM  | C34426   |
| Selenium                                  | 0.038  | 0.011   | 0.050  | J    | mg/L  | 50  | 5/23/2016 6:56:47 PM  | C34426   |
| <b>EPA 200.8: METALS</b>                  |        |         |        |      |       |     | Analyst: DBD          |          |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-20-GW

**Project:** SWMU 10

**Collection Date:** 5/20/2016 10:45:00 AM

**Lab ID:** 1605998-001

**Matrix:** AQUEOUS

**Received Date:** 5/20/2016 3:00:00 PM

| Analyses                               | Result  | MDL      | PQL     | Qual | Units | DF | Date Analyzed        | Batch ID |
|--|---------|----------|---------|------|-------|----|----------------------|----------|
| <b>EPA 200.8: METALS</b>               |         |          |         |      |       |    | Analyst: <b>DBD</b>  |          |
| Antimony                               | ND      | 0.0024   | 0.0050  |      | mg/L  | 5  | 5/25/2016 2:30:09 PM | 25486    |
| Arsenic                                | 0.026   | 0.0029   | 0.010   | *    | mg/L  | 10 | 5/25/2016 3:11:35 PM | 25486    |
| Lead                                   | 0.011   | 0.00084  | 0.0025  |      | mg/L  | 5  | 5/25/2016 2:30:09 PM | 25486    |
| Selenium                               | 0.040   | 0.0021   | 0.010   |      | mg/L  | 10 | 5/25/2016 3:11:35 PM | 25486    |
| <b>EPA METHOD 245.1: MERCURY</b>       |         |          |         |      |       |    | Analyst: <b>pmf</b>  |          |
| Mercury                                | 0.00016 | 0.000053 | 0.00020 | J    | mg/L  | 1  | 5/26/2016 6:31:07 PM | 25512    |
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |         |          |         |      |       |    | Analyst: <b>DAM</b>  |          |
| Acenaphthene                           | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Acenaphthylene                         | ND      | 120      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Aniline                                | ND      | 120      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Anthracene                             | ND      | 120      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Azobenzene                             | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Benz(a)anthracene                      | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Benzo(a)pyrene                         | ND      | 140      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Benzo(b)fluoranthene                   | ND      | 140      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Benzo(g,h,i)perylene                   | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Benzo(k)fluoranthene                   | ND      | 150      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Benzoic acid                           | ND      | 130      | 1000    | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Benzyl alcohol                         | ND      | 150      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Bis(2-chloroethoxy)methane             | ND      | 140      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Bis(2-chloroethyl)ether                | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Bis(2-chloroisopropyl)ether            | ND      | 95       | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Bis(2-ethylhexyl)phthalate             | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| 4-Bromophenyl phenyl ether             | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Butyl benzyl phthalate                 | ND      | 120      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Carbazole                              | ND      | 110      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| 4-Chloro-3-methylphenol                | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| 4-Chloroaniline                        | ND      | 140      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| 2-Chloronaphthalene                    | ND      | 110      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| 2-Chlorophenol                         | ND      | 110      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| 4-Chlorophenyl phenyl ether            | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Chrysene                               | ND      | 140      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Di-n-butyl phthalate                   | ND      | 120      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Di-n-octyl phthalate                   | ND      | 99       | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Dibenz(a,h)anthracene                  | ND      | 130      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| Dibenzofuran                           | ND      | 120      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| 1,2-Dichlorobenzene                    | ND      | 110      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |
| 1,3-Dichlorobenzene                    | ND      | 110      | 500     | D    | µg/L  | 10 | 5/24/2016 4:35:11 PM | 25437    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-20-GW

**Project:** SWMU 10

**Collection Date:** 5/20/2016 10:45:00 AM

**Lab ID:** 1605998-001

**Matrix:** AQUEOUS

**Received Date:** 5/20/2016 3:00:00 PM

| Analyses                               | Result | MDL  | PQL   | Qual | Units | DF  | Date Analyzed         | Batch ID |
|--|--------|------|-------|------|-------|-----|-----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |      |       |      |       |     | Analyst: DAM          |          |
| 1,4-Dichlorobenzene                    | ND     | 120  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 3,3'-Dichlorobenzidine                 | ND     | 120  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Diethyl phthalate                      | ND     | 140  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Dimethyl phthalate                     | ND     | 120  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2,4-Dichlorophenol                     | ND     | 120  | 1000  | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2,4-Dimethylphenol                     | 62000  | 1500 | 5000  |      | µg/L  | 100 | 5/25/2016 11:48:15 AM | 25437    |
| 4,6-Dinitro-2-methylphenol             | ND     | 90   | 1000  | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2,4-Dinitrophenol                      | ND     | 140  | 1000  | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2,4-Dinitrotoluene                     | ND     | 160  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2,6-Dinitrotoluene                     | ND     | 140  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Fluoranthene                           | ND     | 130  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Fluorene                               | ND     | 140  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Hexachlorobenzene                      | ND     | 130  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Hexachlorobutadiene                    | ND     | 110  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Hexachlorocyclopentadiene              | ND     | 110  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Hexachloroethane                       | ND     | 120  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Indeno(1,2,3-cd)pyrene                 | ND     | 150  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Isophorone                             | ND     | 130  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 1-Methylnaphthalene                    | 170    | 150  | 500   | JD   | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2-Methylnaphthalene                    | 240    | 140  | 500   | JD   | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2-Methylphenol                         | 130000 | 2500 | 10000 |      | µg/L  | 200 | 5/25/2016 11:18:11 AM | 25437    |
| 3+4-Methylphenol                       | 190000 | 2300 | 10000 |      | µg/L  | 200 | 5/25/2016 11:18:11 AM | 25437    |
| N-Nitrosodi-n-propylamine              | ND     | 120  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| N-Nitrosodimethylamine                 | ND     | 110  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| N-Nitrosodiphenylamine                 | ND     | 120  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Naphthalene                            | 240    | 130  | 500   | JD   | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2-Nitroaniline                         | ND     | 140  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 3-Nitroaniline                         | ND     | 150  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 4-Nitroaniline                         | ND     | 130  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Nitrobenzene                           | ND     | 140  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2-Nitrophenol                          | ND     | 120  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 4-Nitrophenol                          | ND     | 130  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Pentachlorophenol                      | ND     | 120  | 1000  | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Phenanthrene                           | ND     | 130  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Phenol                                 | 89000  | 990  | 5000  |      | µg/L  | 100 | 5/25/2016 11:48:15 AM | 25437    |
| Pyrene                                 | ND     | 150  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Pyridine                               | ND     | 110  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 1,2,4-Trichlorobenzene                 | ND     | 130  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| 2,4,5-Trichlorophenol                  | ND     | 110  | 500   | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |   |   |
|--------------------|---|---|
| <b>Qualifiers:</b> | * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
|                    | D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
|                    | H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
|                    | ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
|                    | R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
|                    | S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |



# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-20-GW

Project: SWMU 10

Collection Date: 5/20/2016 10:45:00 AM

Lab ID: 1605998-001

Matrix: AQUEOUS

Received Date: 5/20/2016 3:00:00 PM

| Analyses                               | Result | MDL  | PQL      | Qual | Units | DF  | Date Analyzed         | Batch ID |
|--|--------|------|----------|------|-------|-----|-----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |      |          |      |       |     | Analyst: DAM          |          |
| 2,4,6-Trichlorophenol                  | ND     | 120  | 500      | D    | µg/L  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Surr: 2-Fluorophenol                   | 67.7   | 0    | 15-123   | D    | %Rec  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Surr: Phenol-d5                        | 42.0   | 0    | 4.13-124 | D    | %Rec  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Surr: 2,4,6-Tribromophenol             | 92.8   | 0    | 18.4-134 | D    | %Rec  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Surr: Nitrobenzene-d5                  | 107    | 0    | 28.8-134 | D    | %Rec  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Surr: 2-Fluorobiphenyl                 | 96.4   | 0    | 35.9-125 | D    | %Rec  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| Surr: 4-Terphenyl-d14                  | 76.6   | 0    | 15-146   | D    | %Rec  | 10  | 5/24/2016 4:35:11 PM  | 25437    |
| <b>EPA METHOD 8260B: VOLATILES</b>     |        |      |          |      |       |     | Analyst: DJF          |          |
| Benzene                                | 1600   | 9.6  | 100      |      | µg/L  | 100 | 5/26/2016 4:43:13 PM  | A34513   |
| Toluene                                | 4000   | 12   | 100      |      | µg/L  | 100 | 5/26/2016 4:43:13 PM  | A34513   |
| Ethylbenzene                           | 290    | 1.1  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Methyl tert-butyl ether (MTBE)         | 3.8    | 2.1  | 10       | J    | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 1,2,4-Trimethylbenzene                 | 220    | 1.1  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 1,3,5-Trimethylbenzene                 | 64     | 1.2  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 1,2-Dichloroethane (EDC)               | ND     | 1.2  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 1,2-Dibromoethane (EDB)                | ND     | 1.1  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Naphthalene                            | 310    | 0.93 | 20       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 1-Methylnaphthalene                    | 140    | 2.0  | 40       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 2-Methylnaphthalene                    | 210    | 1.6  | 40       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Acetone                                | 1400   | 49   | 100      |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Bromobenzene                           | ND     | 0.98 | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Bromodichloromethane                   | ND     | 1.4  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Bromoform                              | ND     | 1.0  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Bromomethane                           | ND     | 7.8  | 30       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 2-Butanone                             | 630    | 7.4  | 100      |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Carbon disulfide                       | ND     | 6.0  | 100      |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Carbon Tetrachloride                   | ND     | 1.1  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Chlorobenzene                          | ND     | 1.1  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Chloroethane                           | ND     | 1.9  | 20       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Chloroform                             | ND     | 0.89 | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Chloromethane                          | ND     | 2.1  | 30       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 2-Chlorotoluene                        | ND     | 4.0  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 4-Chlorotoluene                        | ND     | 1.3  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| cis-1,2-DCE                            | ND     | 1.2  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| cis-1,3-Dichloropropene                | ND     | 1.1  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 1,2-Dibromo-3-chloropropane            | ND     | 2.3  | 20       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Dibromochloromethane                   | ND     | 0.87 | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| Dibromomethane                         | ND     | 1.2  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |
| 1,2-Dichlorobenzene                    | ND     | 4.0  | 10       |      | µg/L  | 10  | 5/26/2016 12:37:51 PM | A34513   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|             |    |   |    |   |
|-------------|----|---|----|---|
| Qualifiers: | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|             | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|             | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|             | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|             | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|             | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-20-GW

Project: SWMU 10

Collection Date: 5/20/2016 10:45:00 AM

Lab ID: 1605998-001

Matrix: AQUEOUS

Received Date: 5/20/2016 3:00:00 PM

| Analyses                           | Result | MDL  | PQL    | Qual | Units | DF | Date Analyzed         | Batch ID |
|------------------------------------|--------|------|--------|------|-------|----|-----------------------|----------|
| <b>EPA METHOD 8260B: VOLATILES</b> |        |      |        |      |       |    | Analyst: DJF          |          |
| 1,3-Dichlorobenzene                | ND     | 1.4  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,4-Dichlorobenzene                | ND     | 1.4  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Dichlorodifluoromethane            | ND     | 3.6  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,1-Dichloroethane                 | 6.1    | 1.1  | 10     | J    | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,1-Dichloroethene                 | ND     | 1.1  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,2-Dichloropropane                | ND     | 1.1  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,3-Dichloropropane                | ND     | 1.6  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 2,2-Dichloropropane                | ND     | 1.7  | 20     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,1-Dichloropropene                | ND     | 1.3  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Hexachlorobutadiene                | ND     | 2.0  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 2-Hexanone                         | 69     | 8.4  | 100    | J    | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Isopropylbenzene                   | 20     | 1.0  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 4-Isopropyltoluene                 | 2.5    | 1.4  | 10     | J    | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 4-Methyl-2-pentanone               | 63     | 4.3  | 100    | J    | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Methylene Chloride                 | 3.4    | 1.9  | 30     | J    | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| n-Butylbenzene                     | 4.0    | 1.6  | 30     | J    | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| n-Propylbenzene                    | 30     | 1.3  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| sec-Butylbenzene                   | 1.7    | 1.2  | 10     | J    | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Styrene                            | ND     | 1.1  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| tert-Butylbenzene                  | ND     | 1.2  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,1,1,2-Tetrachloroethane          | ND     | 1.1  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,1,2,2-Tetrachloroethane          | ND     | 1.3  | 20     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Tetrachloroethene (PCE)            | ND     | 1.5  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| trans-1,2-DCE                      | ND     | 4.0  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| trans-1,3-Dichloropropene          | ND     | 1.0  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,2,3-Trichlorobenzene             | ND     | 1.1  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,2,4-Trichlorobenzene             | ND     | 1.3  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,1,1-Trichloroethane              | ND     | 0.91 | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,1,2-Trichloroethane              | ND     | 1.3  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Trichloroethene (TCE)              | ND     | 1.8  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Trichlorofluoromethane             | ND     | 2.0  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| 1,2,3-Trichloropropane             | ND     | 2.0  | 20     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Vinyl chloride                     | ND     | 2.0  | 10     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Xylenes, Total                     | 1900   | 3.7  | 15     |      | µg/L  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Surr: 1,2-Dichloroethane-d4        | 96.2   | 0    | 70-130 |      | %Rec  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Surr: 4-Bromofluorobenzene         | 109    | 0    | 70-130 |      | %Rec  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Surr: Dibromofluoromethane         | 100    | 0    | 70-130 |      | %Rec  | 10 | 5/26/2016 12:37:51 PM | A34513   |
| Surr: Toluene-d8                   | 102    | 0    | 70-130 |      | %Rec  | 10 | 5/26/2016 12:37:51 PM | A34513   |

## EPA METHOD 8015D: GASOLINE RANGE

Analyst: DJF

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |   |   |
|--------------------|---|---|
| <b>Qualifiers:</b> | * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
|                    | D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
|                    | H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
|                    | ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
|                    | R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
|                    | S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-20-GW

**Project:** SWMU 10

**Collection Date:** 5/20/2016 10:45:00 AM

**Lab ID:** 1605998-001

**Matrix:** AQUEOUS

**Received Date:** 5/20/2016 3:00:00 PM

| Analyses                                | Result | MDL  | PQL    | Qual | Units | DF | Date Analyzed         | Batch ID |
|---|--------|------|--------|------|-------|----|-----------------------|----------|
| <b>EPA METHOD 8015D: GASOLINE RANGE</b> |        |      |        |      |       |    | Analyst: <b>DJF</b>   |          |
| Gasoline Range Organics (GRO)           | 25     | 0.15 | 0.50   |      | mg/L  | 10 | 5/26/2016 12:37:51 PM | C34513   |
| Surr: BFB                               | 101    | 0    | 70-130 |      | %Rec  | 10 | 5/26/2016 12:37:51 PM | C34513   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

CLIENT: Western Refining Company

Client Sample ID: Trip Blank

Project: SWMU 10

Collection Date:

Lab ID: 1605998-002

Matrix: TRIP BLANK

Received Date: 5/20/2016 3:00:00 PM

| Analyses                           | Result | MDL   | PQL | Qual | Units | DF | Date Analyzed        | Batch ID |
|------------------------------------|--------|-------|-----|------|-------|----|----------------------|----------|
| <b>EPA METHOD 8260B: VOLATILES</b> |        |       |     |      |       |    | Analyst: DJF         |          |
| Benzene                            | ND     | 0.096 | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Toluene                            | ND     | 0.12  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Ethylbenzene                       | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Methyl tert-butyl ether (MTBE)     | ND     | 0.21  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,2,4-Trimethylbenzene             | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,3,5-Trimethylbenzene             | ND     | 0.12  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,2-Dichloroethane (EDC)           | ND     | 0.12  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,2-Dibromoethane (EDB)            | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Naphthalene                        | ND     | 0.093 | 2.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1-Methylnaphthalene                | 1.1    | 0.20  | 4.0 | J    | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 2-Methylnaphthalene                | 1.3    | 0.16  | 4.0 | J    | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Acetone                            | ND     | 4.9   | 10  |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Bromobenzene                       | ND     | 0.098 | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Bromodichloromethane               | ND     | 0.14  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Bromoform                          | ND     | 0.10  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Bromomethane                       | ND     | 0.78  | 3.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 2-Butanone                         | ND     | 0.74  | 10  |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Carbon disulfide                   | ND     | 0.60  | 10  |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Carbon Tetrachloride               | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Chlorobenzene                      | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Chloroethane                       | ND     | 0.19  | 2.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Chloroform                         | ND     | 0.089 | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Chloromethane                      | ND     | 0.21  | 3.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 2-Chlorotoluene                    | ND     | 0.40  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 4-Chlorotoluene                    | ND     | 0.13  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| cis-1,2-DCE                        | ND     | 0.12  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| cis-1,3-Dichloropropene            | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,2-Dibromo-3-chloropropane        | ND     | 0.23  | 2.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Dibromochloromethane               | ND     | 0.087 | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Dibromomethane                     | ND     | 0.12  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,2-Dichlorobenzene                | ND     | 0.40  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,3-Dichlorobenzene                | ND     | 0.14  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,4-Dichlorobenzene                | ND     | 0.14  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Dichlorodifluoromethane            | ND     | 0.36  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,1-Dichloroethane                 | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,1-Dichloroethene                 | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,2-Dichloropropane                | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,3-Dichloropropane                | ND     | 0.16  | 1.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 2,2-Dichloropropane                | ND     | 0.17  | 2.0 |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

CLIENT: Western Refining Company

Client Sample ID: Trip Blank

Project: SWMU 10

Collection Date:

Lab ID: 1605998-002

Matrix: TRIP BLANK

Received Date: 5/20/2016 3:00:00 PM

| Analyses                           | Result | MDL   | PQL    | Qual | Units | DF | Date Analyzed        | Batch ID |
|------------------------------------|--------|-------|--------|------|-------|----|----------------------|----------|
| <b>EPA METHOD 8260B: VOLATILES</b> |        |       |        |      |       |    | Analyst: DJF         |          |
| 1,1-Dichloropropene                | ND     | 0.13  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Hexachlorobutadiene                | ND     | 0.20  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 2-Hexanone                         | ND     | 0.84  | 10     |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Isopropylbenzene                   | ND     | 0.10  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 4-Isopropyltoluene                 | ND     | 0.14  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 4-Methyl-2-pentanone               | ND     | 0.43  | 10     |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Methylene Chloride                 | 0.33   | 0.19  | 3.0    | J    | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| n-Butylbenzene                     | ND     | 0.16  | 3.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| n-Propylbenzene                    | ND     | 0.13  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| sec-Butylbenzene                   | ND     | 0.12  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Styrene                            | ND     | 0.11  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| tert-Butylbenzene                  | ND     | 0.12  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,1,1,2-Tetrachloroethane          | ND     | 0.11  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,1,2,2-Tetrachloroethane          | ND     | 0.13  | 2.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Tetrachloroethene (PCE)            | ND     | 0.15  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| trans-1,2-DCE                      | ND     | 0.40  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| trans-1,3-Dichloropropene          | ND     | 0.10  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,2,3-Trichlorobenzene             | ND     | 0.11  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,2,4-Trichlorobenzene             | ND     | 0.13  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,1,1-Trichloroethane              | ND     | 0.091 | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,1,2-Trichloroethane              | ND     | 0.13  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Trichloroethene (TCE)              | ND     | 0.18  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Trichlorofluoromethane             | ND     | 0.20  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| 1,2,3-Trichloropropane             | ND     | 0.20  | 2.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Vinyl chloride                     | ND     | 0.20  | 1.0    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Xylenes, Total                     | ND     | 0.37  | 1.5    |      | µg/L  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Surr: 1,2-Dichloroethane-d4        | 94.9   | 0     | 70-130 |      | %Rec  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Surr: 4-Bromofluorobenzene         | 111    | 0     | 70-130 |      | %Rec  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Surr: Dibromofluoromethane         | 101    | 0     | 70-130 |      | %Rec  | 1  | 5/26/2016 2:09:48 PM | A34513   |
| Surr: Toluene-d8                   | 103    | 0     | 70-130 |      | %Rec  | 1  | 5/26/2016 2:09:48 PM | A34513   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-21-GW

**Project:** SWMU 10

**Collection Date:** 5/20/2016 9:36:00 AM

**Lab ID:** 1605998-003

**Matrix:** AQUEOUS

**Received Date:** 5/20/2016 3:00:00 PM

| Analyses                               | Result | MDL | PQL | Qual | Units | DF | Date Analyzed        | Batch ID |
|--|--------|-----|-----|------|-------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |     |     |      |       |    | Analyst: <b>DAM</b>  |          |
| Acenaphthene                           | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Acenaphthylene                         | ND     | 12  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Aniline                                | ND     | 12  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Anthracene                             | ND     | 12  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Azobenzene                             | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Benz(a)anthracene                      | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Benzo(a)pyrene                         | ND     | 14  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Benzo(b)fluoranthene                   | ND     | 14  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Benzo(g,h,i)perylene                   | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Benzo(k)fluoranthene                   | ND     | 15  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Benzoic acid                           | ND     | 13  | 100 | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Benzyl alcohol                         | ND     | 15  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Bis(2-chloroethoxy)methane             | ND     | 14  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Bis(2-chloroethyl)ether                | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Bis(2-chloroisopropyl)ether            | ND     | 9.5 | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Bis(2-ethylhexyl)phthalate             | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 4-Bromophenyl phenyl ether             | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Butyl benzyl phthalate                 | ND     | 12  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Carbazole                              | ND     | 11  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 4-Chloro-3-methylphenol                | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 4-Chloroaniline                        | ND     | 14  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2-Chloronaphthalene                    | ND     | 11  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2-Chlorophenol                         | ND     | 11  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 4-Chlorophenyl phenyl ether            | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Chrysene                               | ND     | 14  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Di-n-butyl phthalate                   | ND     | 12  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Di-n-octyl phthalate                   | ND     | 9.9 | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Dibenz(a,h)anthracene                  | ND     | 13  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Dibenzofuran                           | ND     | 12  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 1,2-Dichlorobenzene                    | ND     | 11  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 1,3-Dichlorobenzene                    | ND     | 11  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 1,4-Dichlorobenzene                    | ND     | 12  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 3,3'-Dichlorobenzidine                 | ND     | 12  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Diethyl phthalate                      | ND     | 14  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Dimethyl phthalate                     | ND     | 12  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2,4-Dichlorophenol                     | ND     | 12  | 100 | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2,4-Dimethylphenol                     | ND     | 15  | 50  | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 4,6-Dinitro-2-methylphenol             | ND     | 9.0 | 100 | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2,4-Dinitrophenol                      | ND     | 14  | 100 | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# Hall Environmental Analysis Laboratory, Inc.

## Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

**CLIENT:** Western Refining Company

**Client Sample ID:** SWMU 10-21-GW

**Project:** SWMU 10

**Collection Date:** 5/20/2016 9:36:00 AM

**Lab ID:** 1605998-003

**Matrix:** AQUEOUS

**Received Date:** 5/20/2016 3:00:00 PM

| Analyses                               | Result | MDL | PQL      | Qual | Units | DF | Date Analyzed        | Batch ID |
|--|--------|-----|----------|------|-------|----|----------------------|----------|
| <b>EPA METHOD 8270C: SEMIVOLATILES</b> |        |     |          |      |       |    | Analyst: <b>DAM</b>  |          |
| 2,4-Dinitrotoluene                     | ND     | 16  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2,6-Dinitrotoluene                     | ND     | 14  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Fluoranthene                           | ND     | 13  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Fluorene                               | ND     | 14  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Hexachlorobenzene                      | ND     | 13  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Hexachlorobutadiene                    | ND     | 11  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Hexachlorocyclopentadiene              | ND     | 11  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Hexachloroethane                       | ND     | 12  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Indeno(1,2,3-cd)pyrene                 | ND     | 15  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Isophorone                             | ND     | 13  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 1-Methylnaphthalene                    | ND     | 15  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2-Methylnaphthalene                    | ND     | 14  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2-Methylphenol                         | ND     | 13  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 3+4-Methylphenol                       | ND     | 11  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| N-Nitrosodi-n-propylamine              | ND     | 12  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| N-Nitrosodimethylamine                 | ND     | 11  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| N-Nitrosodiphenylamine                 | ND     | 12  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Naphthalene                            | ND     | 13  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2-Nitroaniline                         | ND     | 14  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 3-Nitroaniline                         | ND     | 15  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 4-Nitroaniline                         | ND     | 13  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Nitrobenzene                           | ND     | 14  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2-Nitrophenol                          | ND     | 12  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 4-Nitrophenol                          | ND     | 13  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Pentachlorophenol                      | ND     | 12  | 100      | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Phenanthrene                           | ND     | 13  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Phenol                                 | ND     | 9.9 | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Pyrene                                 | ND     | 15  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Pyridine                               | ND     | 11  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 1,2,4-Trichlorobenzene                 | ND     | 13  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2,4,5-Trichlorophenol                  | ND     | 11  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| 2,4,6-Trichlorophenol                  | ND     | 12  | 50       | D    | µg/L  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Surr: 2-Fluorophenol                   | 55.1   | 0   | 15-123   | D    | %Rec  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Surr: Phenol-d5                        | 41.1   | 0   | 4.13-124 | D    | %Rec  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Surr: 2,4,6-Tribromophenol             | 70.0   | 0   | 18.4-134 | D    | %Rec  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Surr: Nitrobenzene-d5                  | 74.5   | 0   | 28.8-134 | D    | %Rec  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Surr: 2-Fluorobiphenyl                 | 71.6   | 0   | 35.9-125 | D    | %Rec  | 1  | 5/24/2016 4:04:33 PM | 25437    |
| Surr: 4-Terphenyl-d14                  | 37.0   | 0   | 15-146   | D    | %Rec  | 1  | 5/24/2016 4:04:33 PM | 25437    |

## EPA METHOD 8260B: VOLATILES

Analyst: **DJF**

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |   |   |
|--------------------|---|---|
| <b>Qualifiers:</b> | * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
|                    | D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
|                    | H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
|                    | ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
|                    | R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
|                    | S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |

## Hall Environmental Analysis Laboratory, Inc.

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21-GW

Project: SWMU 10

Collection Date: 5/20/2016 9:36:00 AM

Lab ID: 1605998-003

Matrix: AQUEOUS

Received Date: 5/20/2016 3:00:00 PM

| Analyses                           | Result | MDL   | PQL | Qual | Units | DF | Date Analyzed        | Batch ID |
|------------------------------------|--------|-------|-----|------|-------|----|----------------------|----------|
| <b>EPA METHOD 8260B: VOLATILES</b> |        |       |     |      |       |    | Analyst: DJF         |          |
| Benzene                            | 0.15   | 0.096 | 1.0 | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Toluene                            | 0.30   | 0.12  | 1.0 | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Ethylbenzene                       | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Methyl tert-butyl ether (MTBE)     | 0.80   | 0.21  | 1.0 | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,2,4-Trimethylbenzene             | 0.50   | 0.11  | 1.0 | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,3,5-Trimethylbenzene             | ND     | 0.12  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,2-Dichloroethane (EDC)           | ND     | 0.12  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,2-Dibromoethane (EDB)            | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Naphthalene                        | 1.3    | 0.093 | 2.0 | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1-Methylnaphthalene                | 3.9    | 0.20  | 4.0 | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 2-Methylnaphthalene                | 4.9    | 0.16  | 4.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Acetone                            | 20     | 4.9   | 10  |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Bromobenzene                       | ND     | 0.098 | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Bromodichloromethane               | 1.8    | 0.14  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Bromoform                          | 0.16   | 0.10  | 1.0 | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Bromomethane                       | ND     | 0.78  | 3.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 2-Butanone                         | 5.7    | 0.74  | 10  | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Carbon disulfide                   | ND     | 0.60  | 10  |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Carbon Tetrachloride               | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Chlorobenzene                      | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Chloroethane                       | ND     | 0.19  | 2.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Chloroform                         | 3.4    | 0.089 | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Chloromethane                      | ND     | 0.21  | 3.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 2-Chlorotoluene                    | ND     | 0.40  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 4-Chlorotoluene                    | ND     | 0.13  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| cis-1,2-DCE                        | ND     | 0.12  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| cis-1,3-Dichloropropene            | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,2-Dibromo-3-chloropropane        | ND     | 0.23  | 2.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Dibromochloromethane               | 1.3    | 0.087 | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Dibromomethane                     | ND     | 0.12  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,2-Dichlorobenzene                | ND     | 0.40  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,3-Dichlorobenzene                | ND     | 0.14  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,4-Dichlorobenzene                | ND     | 0.14  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Dichlorodifluoromethane            | ND     | 0.36  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,1-Dichloroethane                 | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,1-Dichloroethene                 | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,2-Dichloropropane                | ND     | 0.11  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,3-Dichloropropane                | ND     | 0.16  | 1.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 2,2-Dichloropropane                | ND     | 0.17  | 2.0 |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |    |   |    |   |
|--------------------|----|---|----|---|
| <b>Qualifiers:</b> | *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
|                    | D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
|                    | H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
|                    | ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
|                    | R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
|                    | S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |



# Hall Environmental Analysis Laboratory, Inc.

Analytical Report

Lab Order 1605998

Date Reported: 6/30/2016

CLIENT: Western Refining Company

Client Sample ID: SWMU 10-21-GW

Project: SWMU 10

Collection Date: 5/20/2016 9:36:00 AM

Lab ID: 1605998-003

Matrix: AQUEOUS

Received Date: 5/20/2016 3:00:00 PM

| Analyses                                | Result | MDL   | PQL    | Qual | Units | DF | Date Analyzed        | Batch ID |
|---|--------|-------|--------|------|-------|----|----------------------|----------|
| <b>EPA METHOD 8260B: VOLATILES</b>      |        |       |        |      |       |    | Analyst: DJF         |          |
| 1,1-Dichloropropene                     | ND     | 0.13  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Hexachlorobutadiene                     | ND     | 0.20  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 2-Hexanone                              | 3.2    | 0.84  | 10     | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Isopropylbenzene                        | ND     | 0.10  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 4-Isopropyltoluene                      | ND     | 0.14  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 4-Methyl-2-pentanone                    | 2.4    | 0.43  | 10     | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Methylene Chloride                      | ND     | 0.19  | 3.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| n-Butylbenzene                          | ND     | 0.16  | 3.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| n-Propylbenzene                         | ND     | 0.13  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| sec-Butylbenzene                        | ND     | 0.12  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Styrene                                 | ND     | 0.11  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| tert-Butylbenzene                       | ND     | 0.12  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,1,1,2-Tetrachloroethane               | ND     | 0.11  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,1,2,2-Tetrachloroethane               | ND     | 0.13  | 2.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Tetrachloroethene (PCE)                 | ND     | 0.15  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| trans-1,2-DCE                           | ND     | 0.40  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| trans-1,3-Dichloropropene               | ND     | 0.10  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,2,3-Trichlorobenzene                  | ND     | 0.11  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,2,4-Trichlorobenzene                  | ND     | 0.13  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,1,1-Trichloroethane                   | ND     | 0.091 | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,1,2-Trichloroethane                   | ND     | 0.13  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Trichloroethene (TCE)                   | ND     | 0.18  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Trichlorofluoromethane                  | ND     | 0.20  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| 1,2,3-Trichloropropane                  | ND     | 0.20  | 2.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Vinyl chloride                          | ND     | 0.20  | 1.0    |      | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Xylenes, Total                          | 0.72   | 0.37  | 1.5    | J    | µg/L  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Surr: 1,2-Dichloroethane-d4             | 102    | 0     | 70-130 |      | %Rec  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Surr: 4-Bromofluorobenzene              | 113    | 0     | 70-130 |      | %Rec  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Surr: Dibromofluoromethane              | 106    | 0     | 70-130 |      | %Rec  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| Surr: Toluene-d8                        | 101    | 0     | 70-130 |      | %Rec  | 1  | 5/26/2016 1:41:29 PM | A34513   |
| <b>EPA METHOD 8015D: GASOLINE RANGE</b> |        |       |        |      |       |    | Analyst: DJF         |          |
| Gasoline Range Organics (GRO)           | 0.047  | 0.015 | 0.050  | J    | mg/L  | 1  | 5/26/2016 1:41:29 PM | C34513   |
| Surr: BFB                               | 102    | 0     | 70-130 |      | %Rec  | 1  | 5/26/2016 1:41:29 PM | C34513   |

Refer to the QC Summary report and sample login checklist for flagged QC data and preservation information.

|                    |   |   |
|--------------------|---|---|
| <b>Qualifiers:</b> | * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
|                    | D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
|                    | H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
|                    | ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
|                    | R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
|                    | S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |

Wet Chemistry by Method 9012B

| Analyte | Result | Qualifier | MDL    | RDL    | Dilution | Analysis date / time | Batch    |
|---------|--------|-----------|--------|--------|----------|----------------------|----------|
| Cyanide | 1.05   |           | 0.0180 | 0.0500 | 10       | 05/31/2016 14:43     | WG876094 |

2

Ic

3

Ss

4

Cn

5

Sr

6

Qc

7

Gl

8

Sc

9

Sc

Wet Chemistry by Method 9012B

L837343-01

Method Blank (MB)

(MB) R3140646-2 05/31/16 13:55

| Analyte | MB Result<br>mg/l | MB Qualifier | MB MDL<br>mg/l | MB RDL<br>mg/l |
|---------|-------------------|--------------|----------------|----------------|
| Cyanide | 0.00218           | 5.0          | 0.00180        | 0.00500        |

L836467-02 Original Sample (OS) • Duplicate (DUP)

(OS) L836467-02 05/31/16 14:01 • (DUP) R3140646-5 05/31/16 14:02

| Analyte | Original Result<br>mg/l | DUP Result<br>mg/l | Dilution | DUP RPD<br>% | DUP Qualifier | DUP RPD Limits<br>% |
|---------|-------------------------|--------------------|----------|--------------|---------------|---------------------|
| Cyanide | 0.00518                 | ND                 | 1        | 10.0         |               | 20                  |

L837343-01 Original Sample (OS) • Duplicate (DUP)

(OS) L837343-01 05/31/16 14:43 • (DUP) R3140646-8 05/31/16 14:44

| Analyte | Original Result<br>mg/l | DUP Result<br>mg/l | Dilution | DUP RPD<br>% | DUP Qualifier | DUP RPD Limits<br>% |
|---------|-------------------------|--------------------|----------|--------------|---------------|---------------------|
| Cyanide | 1.05                    | 1.18               | 10       | 12.0         |               | 20                  |

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCSD)

(LCS) R3140646-3 05/31/16 13:56 • (LCSD) R3140646-4 05/31/16 13:57

| Analyte | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCSD Result<br>mg/l | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|---------|----------------------|--------------------|---------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| Cyanide | 0.100                | 0.103              | 0.103               | 103           | 103            | 90.0-110         |               |                | 0.000    | 20              |

L836467-03 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MSD)

(OS) L836467-03 05/31/16 14:03 • (MS) R3140646-6 05/31/16 14:04 • (MSD) R3140646-7 05/31/16 14:07

| Analyte | Spike Amount<br>mg/l | Original Result<br>mg/l | MS Result<br>mg/l | MSD Result<br>mg/l | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|---------|----------------------|-------------------------|-------------------|--------------------|----------|------------------|--------------|---------------|----------|-----------------|
| Cyanide | 0.200                | 0.00548                 | 0.187             | 0.183              | 1        | 90.0-110         |              | 16            | 2.00     | 20              |

ACCOUNT:

PROJECT:

SDG:

DATE/TIME:

PAGE:



1605998-001I SWMU 10-20-GW

Collected date/time: 05/20/16 10:45

## SAMPLE RESULTS - 01

L839653

ONE LAB. NATIONWIDE.



Wet Chemistry by Method 7199

| Analyte             | Result   | Qualifier | MDL      | RDL      | Dilution | Analysis         | Batch           |
|---------------------|----------|-----------|----------|----------|----------|------------------|-----------------|
|                     | mg/l     |           | mg/l     | mg/l     |          | date / time      |                 |
| Hexavalent Chromium | 0.000500 |           | 0.000150 | 0.000500 | 1        | 05/31/2016 20:03 | <u>WG876252</u> |

<sup>2</sup>Tc<sup>3</sup>Ss<sup>4</sup>Cn<sup>5</sup>Sr<sup>6</sup>Qc<sup>7</sup>Gl<sup>8</sup>Al<sup>9</sup>Sc

ACCOUNT:

Hall Environmental Analysis Laboratory

PROJECT:

SDG:

L839653

DATE/TIME:

06/08/16 10:36



Method Blank (MB)

(MB) R3140841-2 05/31/16 17:59

| Analyte             | MB Result<br>mg/l | MB Qualifier | MB MDL<br>mg/l | MB RDL<br>mg/l |
|---------------------|-------------------|--------------|----------------|----------------|
| Hexavalent Chromium | U                 | 0.000150     | 0.000500       | 0.000500       |

L836719-01 Original Sample (OS) • Duplicate (DUP)

(OS) L836719-01 05/31/16 18:42 • (DUP) R3140841-5 05/31/16 18:53

| Analyte             | Original Result<br>mg/l | DUP Result<br>mg/l | DUP RPD<br>% | DUP Qualifier | DUP RPD Limits<br>% |
|---------------------|-------------------------|--------------------|--------------|---------------|---------------------|
| Hexavalent Chromium | ND                      | ND                 | 0.000        | 1             | 20                  |

Laboratory Control Sample (LCS) • Laboratory Control Sample Duplicate (LCS-D)

(LCS) R3140841-3 05/31/16 18:07 • (LCS-D) R3140841-4 05/31/16 18:15

| Analyte             | Spike Amount<br>mg/l | LCS Result<br>mg/l | LCSD Result<br>mg/l | LCS Rec.<br>% | LCSD Rec.<br>% | Rec. Limits<br>% | LCS Qualifier | LCSD Qualifier | RPD<br>% | RPD Limits<br>% |
|---------------------|----------------------|--------------------|---------------------|---------------|----------------|------------------|---------------|----------------|----------|-----------------|
| Hexavalent Chromium | 0.00200              | 0.00200            | 0.00200             | 100           | 100            | 90.0-110         | 0.000         | 0.000          | 20       | 20              |

L838184-02 Original Sample (OS) • Matrix Spike (MS) • Matrix Spike Duplicate (MS-D)

(OS) L838184-02 05/31/16 21:17 • (MS) R3140841-6 05/31/16 21:25 • (MS-D) R3140841-7 05/31/16 21:33

| Analyte             | Spike Amount<br>mg/l | Original Result<br>mg/l | MS Result<br>mg/l | MSD Result<br>mg/l | MS Rec.<br>% | MSD Rec.<br>% | Dilution | Rec. Limits<br>% | MS Qualifier | MSD Qualifier | RPD<br>% | RPD Limits<br>% |
|---------------------|----------------------|-------------------------|-------------------|--------------------|--------------|---------------|----------|------------------|--------------|---------------|----------|-----------------|
| Hexavalent Chromium | 0.0500               | ND                      | 0.0522            | 0.0512             | 104          | 102           | 1        | 90.0-110         | 2.00         | 2.00          | 20       | 20              |

2 Tl

3 Ss

4 Cn

5 Sr

6 Qc

7 Gl

8 Al

9 Sc



## Abbreviations and Definitions

|                 |   |
|-----------------|---|
| SDG             | Sample Delivery Group.  |
| MDL             | Method Detection Limit.   |
| RDL             | Reported Detection Limit.   |
| ND              | Not detected at the Reporting Limit (or MDL where applicable).  |
| U               | Not detected at the Reporting Limit (or MDL where applicable).  |
| RPD             | Relative Percent Difference.  |
| Original Sample | The non-spiked sample in the prep batch used to determine the Relative Percent Difference (RPD) from a quality control sample. The Original Sample may not be included within the reported SDG. |
| Rec.            | Recovery.   |

| Qualifier | Description |
|-----------|-------------|
|-----------|-------------|

The remainder of this page intentionally left blank, there are no qualifiers applied to this SDG.

1  
H

2  
Tc

3  
Ss

4  
Cn

5  
Sr

6  
Qc

7  
Gl

8  
Al

9  
Sc

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|            |           |                |           |             |                          |          |           |      |          |      |
|------------|-----------|----------------|-----------|-------------|--------------------------|----------|-----------|------|----------|------|
| Sample ID  | MB-25486  | SampType:      | MBLK      | TestCode:   | EPA Method 200.7: Metals |          |           |      |          |      |
| Client ID: | PBW       | Batch ID:      | 25486     | RunNo:      | 34463                    |          |           |      |          |      |
| Prep Date: | 5/24/2016 | Analysis Date: | 5/25/2016 | SeqNo:      | 1062877                  | Units:   | mg/L      |      |          |      |
| Analyte    | Result    | PQL            | SPK value | SPK Ref Val | %REC                     | LowLimit | HighLimit | %RPD | RPDLimit | Qual |

| Analyte   | Result | PQL    | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
|-----------|--------|--------|-----------|-------------|------|----------|-----------|------|----------|------|
| Barium    | ND     | 0.0020 |           |             |      |          |           |      |          |      |
| Beryllium | ND     | 0.0020 |           |             |      |          |           |      |          |      |
| Cadmium   | ND     | 0.0020 |           |             |      |          |           |      |          |      |
| Chromium  | ND     | 0.0060 |           |             |      |          |           |      |          |      |
| Cobalt    | ND     | 0.0060 |           |             |      |          |           |      |          |      |
| Iron      | ND     | 0.020  |           |             |      |          |           |      |          |      |
| Manganese | ND     | 0.0020 |           |             |      |          |           |      |          |      |
| Nickel    | ND     | 0.010  |           |             |      |          |           |      |          |      |
| Silver    | ND     | 0.0050 |           |             |      |          |           |      |          |      |
| Vanadium  | ND     | 0.050  |           |             |      |          |           |      |          |      |
| Zinc      | ND     | 0.010  |           |             |      |          |           |      |          |      |

|            |           |     |                |             |      |           |                          |      |          |      |  |
|------------|-----------|-----|----------------|-------------|------|-----------|--------------------------|------|----------|------|--|
| Sample ID  | LCS-25486 |     | SampType:      | LCS         |      | TestCode: | EPA Method 200.7: Metals |      |          |      |  |
| Client ID: | LCSW      |     | Batch ID:      | 25486       |      | RunNo:    | 34463                    |      |          |      |  |
| Prep Date: | 5/24/2016 |     | Analysis Date: | 5/25/2016   |      | SeqNo:    | 1062878                  |      | Units:   | mg/L |  |
| Analyte    | Result    | PQL | SPK value      | SPK Ref Val | %REC | LowLimit  | HighLimit                | %RPD | RPDLimit | Qual |  |

| Analyte   | Result | PQL    | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
|-----------|--------|--------|-----------|-------------|------|----------|-----------|------|----------|------|
| Barium    | 0.50   | 0.0020 | 0.5000    | 0           | 99.2 | 85       | 115       |      |          |      |
| Beryllium | 0.50   | 0.0020 | 0.5000    | 0           | 100  | 85       | 115       |      |          |      |
| Cadmium   | 0.51   | 0.0020 | 0.5000    | 0           | 102  | 85       | 115       |      |          |      |
| Chromium  | 0.50   | 0.0060 | 0.5000    | 0           | 101  | 85       | 115       |      |          |      |
| Cobalt    | 0.49   | 0.0060 | 0.5000    | 0           | 97.1 | 85       | 115       |      |          |      |
| Iron      | 0.50   | 0.020  | 0.5000    | 0           | 99.9 | 85       | 115       |      |          |      |
| Manganese | 0.48   | 0.0020 | 0.5000    | 0           | 96.8 | 85       | 115       |      |          |      |
| Nickel    | 0.49   | 0.010  | 0.5000    | 0           | 97.2 | 85       | 115       |      |          |      |
| Silver    | 0.099  | 0.0050 | 0.1000    | 0           | 98.8 | 85       | 115       |      |          |      |
| Vanadium  | 0.51   | 0.050  | 0.5000    | 0           | 101  | 85       | 115       |      |          |      |
| Zinc      | 0.50   | 0.010  | 0.5000    | 0           | 101  | 85       | 115       |      |          |      |

|            |             |     |                |             |      |           |                          |      |          |      |  |
|------------|-------------|-----|----------------|-------------|------|-----------|--------------------------|------|----------|------|--|
| Sample ID  | LLLCS-25486 |     | SampType:      | LCSLL       |      | TestCode: | EPA Method 200.7: Metals |      |          |      |  |
| Client ID: | BatchQC     |     | Batch ID:      | 25486       |      | RunNo:    | 34463                    |      |          |      |  |
| Prep Date: | 5/24/2016   |     | Analysis Date: | 5/25/2016   |      | SeqNo:    | 1062880                  |      | Units:   | mg/L |  |
| Analyte    | Result      | PQL | SPK value      | SPK Ref Val | %REC | LowLimit  | HighLimit                | %RPD | RPDLimit | Qual |  |

| Analyte   | Result | PQL    | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
|-----------|--------|--------|-----------|-------------|------|----------|-----------|------|----------|------|
| Barium    | 0.0013 | 0.0020 | 0.002000  | 0           | 66.5 | 50       | 150       |      |          | J    |
| Beryllium | 0.0020 | 0.0020 | 0.002000  | 0           | 103  | 50       | 150       |      |          |      |
| Cadmium   | 0.0022 | 0.0020 | 0.002000  | 0           | 110  | 50       | 150       |      |          |      |
| Chromium  | 0.0050 | 0.0060 | 0.006000  | 0           | 82.8 | 50       | 150       |      |          | J    |
| Cobalt    | 0.0063 | 0.0060 | 0.006000  | 0           | 105  | 50       | 150       |      |          |      |
| Iron      | 0.024  | 0.020  | 0.02000   | 0           | 121  | 50       | 150       |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

R RPD outside accepted recovery limits

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

P Sample pH Not In Range

RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|            |             |                |           |             |                          |          |           |      |          |      |
|------------|-------------|----------------|-----------|-------------|--------------------------|----------|-----------|------|----------|------|
| Sample ID  | LLLCS-25486 | SampType:      | LCSLL     | TestCode:   | EPA Method 200.7: Metals |          |           |      |          |      |
| Client ID: | BatchQC     | Batch ID:      | 25486     | RunNo:      | 34463                    |          |           |      |          |      |
| Prep Date: | 5/24/2016   | Analysis Date: | 5/25/2016 | SeqNo:      | 1062880                  | Units:   | mg/L      |      |          |      |
| Analyte    | Result      | PQL            | SPK value | SPK Ref Val | %REC                     | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
| Manganese  | 0.0027      | 0.0020         | 0.002000  | 0           | 136                      | 50       | 150       |      |          |      |
| Nickel     | 0.0056      | 0.010          | 0.005000  | 0           | 113                      | 50       | 150       |      |          | J    |
| Silver     | 0.0054      | 0.0050         | 0.005000  | 0           | 109                      | 50       | 150       |      |          |      |
| Vanadium   | 0.0096      | 0.050          | 0.01000   | 0           | 95.8                     | 50       | 150       |      |          | J    |
| Zinc       | 0.0063      | 0.010          | 0.005000  | 0           | 126                      | 50       | 150       |      |          | J    |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified



# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

**Client:** Western Refining Company

**Project:** SWMU 10

|                       |                                 |   |
|-----------------------|---------------------------------|---|
| Sample ID <b>MB-C</b> | SampType: <b>MBLK</b>           | TestCode: <b>EPA Method 200.7: Dissolved Metals</b> |
| Client ID: <b>PBW</b> | Batch ID: <b>C34463</b>         | RunNo: <b>34463</b>                                 |
| Prep Date:            | Analysis Date: <b>5/25/2016</b> | SeqNo: <b>1062977</b> Units: <b>mg/L</b>            |

| Analyte   | Result | PQL    | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
|-----------|--------|--------|-----------|-------------|------|----------|-----------|------|----------|------|
| Barium    | ND     | 0.0020 |           |             |      |          |           |      |          |      |
| Beryllium | ND     | 0.0020 |           |             |      |          |           |      |          |      |
| Cadmium   | ND     | 0.0020 |           |             |      |          |           |      |          |      |
| Chromium  | ND     | 0.0060 |           |             |      |          |           |      |          |      |
| Cobalt    | ND     | 0.0060 |           |             |      |          |           |      |          |      |
| Iron      | ND     | 0.020  |           |             |      |          |           |      |          |      |
| Manganese | ND     | 0.0020 |           |             |      |          |           |      |          |      |
| Nickel    | ND     | 0.010  |           |             |      |          |           |      |          |      |
| Silver    | ND     | 0.0050 |           |             |      |          |           |      |          |      |
| Vanadium  | ND     | 0.050  |           |             |      |          |           |      |          |      |
| Zinc      | ND     | 0.010  |           |             |      |          |           |      |          |      |

|                        |                                 |   |
|------------------------|---------------------------------|---|
| Sample ID <b>LCS-C</b> | SampType: <b>LCS</b>            | TestCode: <b>EPA Method 200.7: Dissolved Metals</b> |
| Client ID: <b>LCSW</b> | Batch ID: <b>C34463</b>         | RunNo: <b>34463</b>                                 |
| Prep Date:             | Analysis Date: <b>5/25/2016</b> | SeqNo: <b>1062978</b> Units: <b>mg/L</b>            |

| Analyte   | Result | PQL    | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
|-----------|--------|--------|-----------|-------------|------|----------|-----------|------|----------|------|
| Barium    | 0.49   | 0.0020 | 0.5000    | 0           | 97.1 | 85       | 115       |      |          |      |
| Beryllium | 0.51   | 0.0020 | 0.5000    | 0           | 101  | 85       | 115       |      |          |      |
| Cadmium   | 0.49   | 0.0020 | 0.5000    | 0           | 98.8 | 85       | 115       |      |          |      |
| Chromium  | 0.48   | 0.0060 | 0.5000    | 0           | 96.7 | 85       | 115       |      |          |      |
| Cobalt    | 0.47   | 0.0060 | 0.5000    | 0           | 93.9 | 85       | 115       |      |          |      |
| Iron      | 0.49   | 0.020  | 0.5000    | 0           | 97.7 | 85       | 115       |      |          |      |
| Manganese | 0.48   | 0.0020 | 0.5000    | 0           | 95.9 | 85       | 115       |      |          |      |
| Nickel    | 0.47   | 0.010  | 0.5000    | 0           | 93.1 | 85       | 115       |      |          |      |
| Silver    | 0.10   | 0.0050 | 0.1000    | 0           | 100  | 85       | 115       |      |          |      |
| Vanadium  | 0.50   | 0.050  | 0.5000    | 0           | 101  | 85       | 115       |      |          |      |
| Zinc      | 0.48   | 0.010  | 0.5000    | 0           | 96.3 | 85       | 115       |      |          |      |

|                           |                                 |   |
|---------------------------|---------------------------------|---|
| Sample ID <b>LLCS-C</b>   | SampType: <b>LCSLL</b>          | TestCode: <b>EPA Method 200.7: Dissolved Metals</b> |
| Client ID: <b>BatchQC</b> | Batch ID: <b>C34463</b>         | RunNo: <b>34463</b>                                 |
| Prep Date:                | Analysis Date: <b>5/25/2016</b> | SeqNo: <b>1062979</b> Units: <b>mg/L</b>            |

| Analyte   | Result | PQL    | SPK value | SPK Ref Val | %REC | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
|-----------|--------|--------|-----------|-------------|------|----------|-----------|------|----------|------|
| Barium    | 0.0022 | 0.0020 | 0.002000  | 0           | 110  | 50       | 150       |      |          |      |
| Beryllium | 0.0022 | 0.0020 | 0.002000  | 0           | 108  | 50       | 150       |      |          |      |
| Cadmium   | 0.0020 | 0.0020 | 0.002000  | 0           | 99.5 | 50       | 150       |      |          | J    |
| Chromium  | 0.0055 | 0.0060 | 0.006000  | 0           | 91.3 | 50       | 150       |      |          | J    |
| Cobalt    | 0.0071 | 0.0060 | 0.006000  | 0           | 118  | 50       | 150       |      |          |      |
| Iron      | 0.026  | 0.020  | 0.02000   | 0           | 128  | 50       | 150       |      |          |      |

### Qualifiers:

|   |   |
|---|---|
| * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
| D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
| H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
| ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
| R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
| S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|            |         |        |                          |             |      |  |           |             |          |      |
|------------|---------|--------|--------------------------|-------------|------|--|-----------|-------------|----------|------|
| Sample ID  | LLLCS-C |        | SampType: LCSLL          |             |      | TestCode: EPA Method 200.7: Dissolved Metals |           |             |          |      |
| Client ID: | BatchQC |        | Batch ID: C34463         |             |      | RunNo: 34463                                 |           |             |          |      |
| Prep Date: |         |        | Analysis Date: 5/25/2016 |             |      | SeqNo: 1062979                               |           | Units: mg/L |          |      |
| Analyte    | Result  | PQL    | SPK value                | SPK Ref Val | %REC | LowLimit                                     | HighLimit | %RPD        | RPDLimit | Qual |
| Manganese  | 0.0021  | 0.0020 | 0.002000                 | 0           | 104  | 50   | 150       |             |          |      |
| Nickel     | 0.0042  | 0.010  | 0.005000                 | 0           | 83.8 | 50   | 150       |             |          | J    |
| Silver     | 0.0055  | 0.0050 | 0.005000                 | 0           | 109  | 50   | 150       |             |          |      |
| Vanadium   | 0.011   | 0.050  | 0.01000                  | 0           | 109  | 50   | 150       |             |          | J    |
| Zinc       | 0.0060  | 0.010  | 0.005000                 | 0           | 121  | 50   | 150       |             |          | J    |

### Qualifiers:

|    |   |    |   |
|----|---|----|---|
| *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
| D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
| H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
| ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
| R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
| S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

**Client:** Western Refining Company

**Project:** SWMU 10

|            |           |         |                |             |      |           |                   |      |          |      |  |
|------------|-----------|---------|----------------|-------------|------|-----------|-------------------|------|----------|------|--|
| Sample ID  | MB-25486  |         | SampType:      | MBLK        |      | TestCode: | EPA 200.8: Metals |      |          |      |  |
| Client ID: | PBW       |         | Batch ID:      | 25486       |      | RunNo:    | 34477             |      |          |      |  |
| Prep Date: | 5/24/2016 |         | Analysis Date: | 5/25/2016   |      | SeqNo:    | 1063236           |      | Units:   | mg/L |  |
| Analyte    | Result    | PQL     | SPK value      | SPK Ref Val | %REC | LowLimit  | HighLimit         | %RPD | RPDLimit | Qual |  |
| Antimony   | ND        | 0.0010  |                |             |      |           |                   |      |          |      |  |
| Arsenic    | ND        | 0.0010  |                |             |      |           |                   |      |          |      |  |
| Lead       | ND        | 0.00050 |                |             |      |           |                   |      |          |      |  |
| Selenium   | ND        | 0.0010  |                |             |      |           |                   |      |          |      |  |

|            |             |         |                          |             |                             |          |             |      |          |      |
|------------|-------------|---------|--------------------------|-------------|-----------------------------|----------|-------------|------|----------|------|
| Sample ID  | MSLCS-25486 |         | SampType: LCS            |             | TestCode: EPA 200.8: Metals |          |             |      |          |      |
| Client ID: | LCSW        |         | Batch ID: 25486          |             | RunNo: 34477                |          |             |      |          |      |
| Prep Date: | 5/24/2016   |         | Analysis Date: 5/25/2016 |             | SeqNo: 1063238              |          | Units: mg/L |      |          |      |
| Analyte    | Result      | PQL     | SPK value                | SPK Ref Val | %REC                        | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Antimony   | 0.026       | 0.0010  | 0.02500                  | 0           | 103                         | 85       | 115         |      |          |      |
| Arsenic    | 0.023       | 0.0010  | 0.02500                  | 0           | 92.3                        | 85       | 115         |      |          |      |
| Lead       | 0.012       | 0.00050 | 0.01250                  | 0           | 99.9                        | 85       | 115         |      |          |      |
| Selenium   | 0.023       | 0.0010  | 0.02500                  | 0           | 93.7                        | 85       | 115         |      |          |      |

|            |              |         |                          |             |                             |          |             |      |          |      |
|------------|--------------|---------|--------------------------|-------------|-----------------------------|----------|-------------|------|----------|------|
| Sample ID  | MSLLCS-25486 |         | SampType: LCSLL          |             | TestCode: EPA 200.8: Metals |          |             |      |          |      |
| Client ID: | BatchQC      |         | Batch ID: 25486          |             | RunNo: 34477                |          |             |      |          |      |
| Prep Date: | 5/24/2016    |         | Analysis Date: 5/25/2016 |             | SeqNo: 1063240              |          | Units: mg/L |      |          |      |
| Analyte    | Result       | PQL     | SPK value                | SPK Ref Val | %REC                        | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Antimony   | 0.00078      | 0.0010  | 0.001000                 | 0           | 78.0                        | 50       | 150         |      |          | J    |
| Arsenic    | 0.00086      | 0.0010  | 0.001000                 | 0           | 86.2                        | 50       | 150         |      |          | J    |
| Lead       | 0.00055      | 0.00050 | 0.0005000                | 0           | 110                         | 50       | 150         |      |          |      |
| Selenium   | 0.00090      | 0.0010  | 0.001000                 | 0           | 90.1                        | 50       | 150         |      |          | J    |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|            |        |         |                          |             |                                       |          |             |      |          |      |
|------------|--------|---------|--------------------------|-------------|---------------------------------------|----------|-------------|------|----------|------|
| Sample ID  | LCS    |         | SampType: LCS            |             | TestCode: EPA 200.8: Dissolved Metals |          |             |      |          |      |
| Client ID: | LCSW   |         | Batch ID: C34426         |             | RunNo: 34426                          |          |             |      |          |      |
| Prep Date: |        |         | Analysis Date: 5/23/2016 |             | SeqNo: 1061595                        |          | Units: mg/L |      |          |      |
| Analyte    | Result | PQL     | SPK value                | SPK Ref Val | %REC                                  | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Antimony   | 0.025  | 0.0010  | 0.02500                  | 0           | 100                                   | 85       | 115         |      |          |      |
| Arsenic    | 0.025  | 0.0010  | 0.02500                  | 0           | 99.1                                  | 85       | 115         |      |          |      |
| Lead       | 0.013  | 0.00050 | 0.01250                  | 0           | 103                                   | 85       | 115         |      |          |      |
| Selenium   | 0.026  | 0.0010  | 0.02500                  | 0           | 103                                   | 85       | 115         |      |          |      |

|            |         |         |                          |             |                                       |          |             |      |          |      |
|------------|---------|---------|--------------------------|-------------|---------------------------------------|----------|-------------|------|----------|------|
| Sample ID  | LCSLL   |         | SampType: LCSLL          |             | TestCode: EPA 200.8: Dissolved Metals |          |             |      |          |      |
| Client ID: | BatchQC |         | Batch ID: C34426         |             | RunNo: 34426                          |          |             |      |          |      |
| Prep Date: |         |         | Analysis Date: 5/23/2016 |             | SeqNo: 1061599                        |          | Units: mg/L |      |          |      |
| Analyte    | Result  | PQL     | SPK value                | SPK Ref Val | %REC                                  | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Antimony   | 0.0010  | 0.0010  | 0.001000                 | 0           | 100                                   | 50       | 150         |      |          |      |
| Arsenic    | 0.00097 | 0.0010  | 0.001000                 | 0           | 96.6                                  | 50       | 150         |      |          | J    |
| Lead       | 0.00053 | 0.00050 | 0.0005000                | 0           | 106                                   | 50       | 150         |      |          |      |
| Selenium   | 0.00097 | 0.0010  | 0.001000                 | 0           | 97.1                                  | 50       | 150         |      |          | J    |

|            |        |                          |           |             |                                       |          |             |      |          |      |
|------------|--------|--------------------------|-----------|-------------|---------------------------------------|----------|-------------|------|----------|------|
| Sample ID  | MB     | SampType: MBLK           |           |             | TestCode: EPA 200.8: Dissolved Metals |          |             |      |          |      |
| Client ID: | PBW    | Batch ID: C34426         |           |             | RunNo: 34426                          |          |             |      |          |      |
| Prep Date: |        | Analysis Date: 5/23/2016 |           |             | SeqNo: 1061601                        |          | Units: mg/L |      |          |      |
| Analyte    | Result | PQL                      | SPK value | SPK Ref Val | %REC                                  | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Antimony   | ND     | 0.0010                   |           |             |                                       |          |             |      |          |      |
| Arsenic    | ND     | 0.0010                   |           |             |                                       |          |             |      |          |      |
| Lead       | ND     | 0.00050                  |           |             |                                       |          |             |      |          |      |
| Selenium   | ND     | 0.0010                   |           |             |                                       |          |             |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|            |           |         |                          |             |                                     |          |             |      |          |      |
|------------|-----------|---------|--------------------------|-------------|-------------------------------------|----------|-------------|------|----------|------|
| Sample ID  | MB-25512  |         | SampType: MBLK           |             | TestCode: EPA Method 245.1: Mercury |          |             |      |          |      |
| Client ID: | PBW       |         | Batch ID: 25512          |             | RunNo: 34519                        |          |             |      |          |      |
| Prep Date: | 5/25/2016 |         | Analysis Date: 5/26/2016 |             | SeqNo: 1064697                      |          | Units: mg/L |      |          |      |
| Analyte    | Result    | PQL     | SPK value                | SPK Ref Val | %REC                                | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Mercurv    | 0.000076  | 0.00020 |                          |             |                                     |          |             |      |          | J    |

|            |           |         |                          |             |                                     |          |             |      |          |      |
|------------|-----------|---------|--------------------------|-------------|-------------------------------------|----------|-------------|------|----------|------|
| Sample ID  | LCS-25512 |         | SampType: LCS            |             | TestCode: EPA Method 245.1: Mercury |          |             |      |          |      |
| Client ID: | LCSW      |         | Batch ID: 25512          |             | RunNo: 34519                        |          |             |      |          |      |
| Prep Date: | 5/25/2016 |         | Analysis Date: 5/26/2016 |             | SeqNo: 1064698                      |          | Units: mg/L |      |          |      |
| Analyte    | Result    | PQL     | SPK value                | SPK Ref Val | %REC                                | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Mercury    | 0.0051    | 0.00020 | 0.005000                 | 0           | 103                                 | 80       | 120         |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|            |            |                |                  |             |                                 |          |             |      |          |      |
|------------|------------|----------------|------------------|-------------|---------------------------------|----------|-------------|------|----------|------|
| Sample ID  | <b>MB</b>  | SampType:      | <b>MBLK</b>      | TestCode:   | <b>EPA Method 300.0: Anions</b> |          |             |      |          |      |
| Client ID: | <b>PBW</b> | Batch ID:      | <b>R34434</b>    | RunNo:      | <b>34434</b>                    |          |             |      |          |      |
| Prep Date: |            | Analysis Date: | <b>5/23/2016</b> | SeqNo:      | <b>1061743</b>                  | Units:   | <b>mg/L</b> |      |          |      |
| Analyte    | Result     | PQL            | SPK value        | SPK Ref Val | %REC                            | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Fluoride   | ND         | 0.10           |                  |             |                                 |          |             |      |          |      |
| Sulfate    | ND         | 0.50           |                  |             |                                 |          |             |      |          |      |

|            |             |                |                  |             |                                 |          |             |      |          |      |
|------------|-------------|----------------|------------------|-------------|---------------------------------|----------|-------------|------|----------|------|
| Sample ID  | <b>LCSb</b> | SampType:      | <b>LCS</b>       | TestCode:   | <b>EPA Method 300.0: Anions</b> |          |             |      |          |      |
| Client ID: | <b>LCSW</b> | Batch ID:      | <b>R34434</b>    | RunNo:      | <b>34434</b>                    |          |             |      |          |      |
| Prep Date: |             | Analysis Date: | <b>5/23/2016</b> | SeqNo:      | <b>1061747</b>                  | Units:   | <b>mg/L</b> |      |          |      |
| Analyte    | Result      | PQL            | SPK value        | SPK Ref Val | %REC                            | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Fluoride   | 0.54        | 0.10           | 0.5000           | 0           | 108                             | 90       | 110         |      |          |      |
| Sulfate    | 9.9         | 0.50           | 10.00            | 0           | 98.5                            | 90       | 110         |      |          |      |

|            |            |                |                  |             |                                 |          |             |      |          |      |
|------------|------------|----------------|------------------|-------------|---------------------------------|----------|-------------|------|----------|------|
| Sample ID  | <b>MB</b>  | SampType:      | <b>MBLK</b>      | TestCode:   | <b>EPA Method 300.0: Anions</b> |          |             |      |          |      |
| Client ID: | <b>PBW</b> | Batch ID:      | <b>R34487</b>    | RunNo:      | <b>34487</b>                    |          |             |      |          |      |
| Prep Date: |            | Analysis Date: | <b>5/25/2016</b> | SeqNo:      | <b>1063631</b>                  | Units:   | <b>mg/L</b> |      |          |      |
| Analyte    | Result     | PQL            | SPK value        | SPK Ref Val | %REC                            | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Chloride   | ND         | 0.50           |                  |             |                                 |          |             |      |          |      |

|            |             |                |                  |             |                                 |          |             |      |          |      |
|------------|-------------|----------------|------------------|-------------|---------------------------------|----------|-------------|------|----------|------|
| Sample ID  | <b>LCS</b>  | SampType:      | <b>LCS</b>       | TestCode:   | <b>EPA Method 300.0: Anions</b> |          |             |      |          |      |
| Client ID: | <b>LCSW</b> | Batch ID:      | <b>R34487</b>    | RunNo:      | <b>34487</b>                    |          |             |      |          |      |
| Prep Date: |             | Analysis Date: | <b>5/25/2016</b> | SeqNo:      | <b>1063632</b>                  | Units:   | <b>mg/L</b> |      |          |      |
| Analyte    | Result      | PQL            | SPK value        | SPK Ref Val | %REC                            | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Chloride   | 4.9         | 0.50           | 5.000            | 0           | 98.3                            | 90       | 110         |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

R RPD outside accepted recovery limits

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

P Sample pH Not In Range

RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                                |           |     |                          |             |  |          |             |      |          |      |
|--------------------------------|-----------|-----|--------------------------|-------------|--|----------|-------------|------|----------|------|
| Sample ID                      | MB-25494  |     | SampType: MBLK           |             | TestCode: EPA Method 8015M/D: Diesel Range |          |             |      |          |      |
| Client ID:                     | PBW       |     | Batch ID: 25494          |             | RunNo: 34483                               |          |             |      |          |      |
| Prep Date:                     | 5/25/2016 |     | Analysis Date: 5/25/2016 |             | SeqNo: 1063745                             |          | Units: mg/L |      |          |      |
| Analyte                        | Result    | PQL | SPK value                | SPK Ref Val | %REC                                       | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| Diesel Range Organics (DRO)    | ND        | 1.0 |                          |             |  |          |             |      |          |      |
| Motor Oil Range Organics (MRO) | ND        | 5.0 |                          |             |  |          |             |      |          |      |
| Surr: DNOP                     | 0.98      |     | 1.000                    |             | 97.7                                       | 70       | 141         |      |          |      |

|                             |           |     |                |             |      |           |                                  |      |          |      |  |
|-----------------------------|-----------|-----|----------------|-------------|------|-----------|----------------------------------|------|----------|------|--|
| Sample ID                   | LCS-25494 |     | SampType:      | LCS         |      | TestCode: | EPA Method 8015M/D: Diesel Range |      |          |      |  |
| Client ID:                  | LCSW      |     | Batch ID:      | 25494       |      | RunNo:    | 34489                            |      |          |      |  |
| Prep Date:                  | 5/25/2016 |     | Analysis Date: | 5/26/2016   |      | SeqNo:    | 1064130                          |      | Units:   | mg/L |  |
| Analyte                     | Result    | PQL | SPK value      | SPK Ref Val | %REC | LowLimit  | HighLimit                        | %RPD | RPDLimit | Qual |  |
| Diesel Range Organics (DRO) | 5.6       | 1.0 | 5.000          | 0           | 111  | 71.3      | 139                              |      |          |      |  |
| Surr: DNOP                  | 0.50      |     | 0.5000         |             | 101  | 70        | 141                              |      |          |      |  |

### Qualifiers:

|   |   |
|---|---|
| * Value exceeds Maximum Contaminant Level.              | B Analyte detected in the associated Method Blank           |
| D Sample Diluted Due to Matrix                          | E Value above quantitation range                            |
| H Holding times for preparation or analysis exceeded    | J Analyte detected below quantitation limits                |
| ND Not Detected at the Reporting Limit                  | P Sample pH Not In Range                                    |
| R RPD outside accepted recovery limits                  | RL Reporting Detection Limit                                |
| S % Recovery outside of range due to dilution or matrix | W Sample container temperature is out of limit as specified |

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                                |        |                |           |             |                             |          |           |      |          |      |
|--------------------------------|--------|----------------|-----------|-------------|-----------------------------|----------|-----------|------|----------|------|
| Sample ID                      | rb1    | SampType:      | MBLK      | TestCode:   | EPA Method 8260B: VOLATILES |          |           |      |          |      |
| Client ID:                     | PBW    | Batch ID:      | A34513    | RunNo:      | 34513                       |          |           |      |          |      |
| Prep Date:                     |        | Analysis Date: | 5/26/2016 | SeqNo:      | 1064513                     | Units:   | µg/L      |      |          |      |
| Analyte                        | Result | PQL            | SPK value | SPK Ref Val | %REC                        | LowLimit | HighLimit | %RPD | RPDLimit | Qual |
| Benzene                        | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Toluene                        | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Ethylbenzene                   | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Methyl tert-butyl ether (MTBE) | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,2,4-Trimethylbenzene         | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,3,5-Trimethylbenzene         | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,2-Dichloroethane (EDC)       | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,2-Dibromoethane (EDB)        | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Naphthalene                    | ND     | 2.0            |           |             |                             |          |           |      |          |      |
| 1-Methylnaphthalene            | ND     | 4.0            |           |             |                             |          |           |      |          |      |
| 2-Methylnaphthalene            | ND     | 4.0            |           |             |                             |          |           |      |          |      |
| Acetone                        | ND     | 10             |           |             |                             |          |           |      |          |      |
| Bromobenzene                   | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Bromodichloromethane           | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Bromoform                      | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Bromomethane                   | ND     | 3.0            |           |             |                             |          |           |      |          |      |
| 2-Butanone                     | ND     | 10             |           |             |                             |          |           |      |          |      |
| Carbon disulfide               | ND     | 10             |           |             |                             |          |           |      |          |      |
| Carbon Tetrachloride           | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Chlorobenzene                  | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Chloroethane                   | ND     | 2.0            |           |             |                             |          |           |      |          |      |
| Chloroform                     | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Chloromethane                  | ND     | 3.0            |           |             |                             |          |           |      |          |      |
| 2-Chlorotoluene                | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 4-Chlorotoluene                | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| cis-1,2-DCE                    | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| cis-1,3-Dichloropropene        | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,2-Dibromo-3-chloropropane    | ND     | 2.0            |           |             |                             |          |           |      |          |      |
| Dibromochloromethane           | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Dibromomethane                 | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,2-Dichlorobenzene            | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,3-Dichlorobenzene            | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,4-Dichlorobenzene            | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| Dichlorodifluoromethane        | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,1-Dichloroethane             | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,1-Dichloroethene             | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,2-Dichloropropane            | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 1,3-Dichloropropane            | ND     | 1.0            |           |             |                             |          |           |      |          |      |
| 2,2-Dichloropropane            | ND     | 2.0            |           |             |                             |          |           |      |          |      |

### Qualifiers:

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D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
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W Sample container temperature is out of limit as specified



# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

| Sample ID <b>rb1</b>        | SampType: <b>MBLK</b>           |     |           | TestCode: <b>EPA Method 8260B: VOLATILES</b> |      |                    |           |      |          |      |
|-----------------------------|---------------------------------|-----|-----------|--|------|--------------------|-----------|------|----------|------|
| Client ID: <b>PBW</b>       | Batch ID: <b>A34513</b>         |     |           | RunNo: <b>34513</b>                          |      |                    |           |      |          |      |
| Prep Date:                  | Analysis Date: <b>5/26/2016</b> |     |           | SeqNo: <b>1064513</b>                        |      | Units: <b>µg/L</b> |           |      |          |      |
| Analyte                     | Result                          | PQL | SPK value | SPK Ref Val                                  | %REC | LowLimit           | HighLimit | %RPD | RPDLimit | Qual |
| 1,1-Dichloropropene         | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| Hexachlorobutadiene         | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 2-Hexanone                  | ND                              | 10  |           |  |      |                    |           |      |          |      |
| Isopropylbenzene            | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 4-Isopropyltoluene          | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 4-Methyl-2-pentanone        | ND                              | 10  |           |  |      |                    |           |      |          |      |
| Methylene Chloride          | 0.45                            | 3.0 |           |  |      |                    |           |      |          | J    |
| n-Butylbenzene              | ND                              | 3.0 |           |  |      |                    |           |      |          |      |
| n-Propylbenzene             | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| sec-Butylbenzene            | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| Styrene                     | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| tert-Butylbenzene           | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 1,1,1,2-Tetrachloroethane   | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 1,1,2,2-Tetrachloroethane   | ND                              | 2.0 |           |  |      |                    |           |      |          |      |
| Tetrachloroethene (PCE)     | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| trans-1,2-DCE               | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| trans-1,3-Dichloropropene   | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 1,2,3-Trichlorobenzene      | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 1,2,4-Trichlorobenzene      | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 1,1,1-Trichloroethane       | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 1,1,2-Trichloroethane       | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| Trichloroethene (TCE)       | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| Trichlorofluoromethane      | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| 1,2,3-Trichloropropane      | ND                              | 2.0 |           |  |      |                    |           |      |          |      |
| Vinyl chloride              | ND                              | 1.0 |           |  |      |                    |           |      |          |      |
| Xylenes, Total              | ND                              | 1.5 |           |  |      |                    |           |      |          |      |
| Surr: 1,2-Dichloroethane-d4 | 9.6                             |     | 10.00     |  | 96.2 | 70                 | 130       |      |          |      |
| Surr: 4-Bromofluorobenzene  | 11                              |     | 10.00     |  | 109  | 70                 | 130       |      |          |      |
| Surr: Dibromofluoromethane  | 10                              |     | 10.00     |  | 101  | 70                 | 130       |      |          |      |
| Surr: Toluene-d8            | 10                              |     | 10.00     |  | 101  | 70                 | 130       |      |          |      |

| Sample ID <b>100ng lcs</b> | SampType: <b>LCS</b>            |     |           | TestCode: <b>EPA Method 8260B: VOLATILES</b> |      |                    |           |      |          |      |
|----------------------------|---------------------------------|-----|-----------|--|------|--------------------|-----------|------|----------|------|
| Client ID: <b>LCSW</b>     | Batch ID: <b>A34513</b>         |     |           | RunNo: <b>34513</b>                          |      |                    |           |      |          |      |
| Prep Date:                 | Analysis Date: <b>5/26/2016</b> |     |           | SeqNo: <b>1064514</b>                        |      | Units: <b>µg/L</b> |           |      |          |      |
| Analyte                    | Result                          | PQL | SPK value | SPK Ref Val                                  | %REC | LowLimit           | HighLimit | %RPD | RPDLimit | Qual |
| Benzene                    | 19                              | 1.0 | 20.00     | 0  | 95.7 | 70                 | 130       |      |          |      |
| Toluene                    | 20                              | 1.0 | 20.00     | 0  | 101  | 70                 | 130       |      |          |      |
| Chlorobenzene              | 20                              | 1.0 | 20.00     | 0  | 99.4 | 70                 | 130       |      |          |      |

### Qualifiers:

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S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

P Sample pH Not In Range

RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                             |           |                          |           |             |                                       |          |             |      |          |      |
|-----------------------------|-----------|--------------------------|-----------|-------------|---------------------------------------|----------|-------------|------|----------|------|
| Sample ID                   | 100ng lcs | SampType: LCS            |           |             | TestCode: EPA Method 8260B: VOLATILES |          |             |      |          |      |
| Client ID:                  | LCSW      | Batch ID: A34513         |           |             | RunNo: 34513                          |          |             |      |          |      |
| Prep Date:                  |           | Analysis Date: 5/26/2016 |           |             | SeqNo: 1064514                        |          | Units: µg/L |      |          |      |
| Analyte                     | Result    | PQL                      | SPK value | SPK Ref Val | %REC                                  | LowLimit | HighLimit   | %RPD | RPDLimit | Qual |
| 1,1-Dichloroethene          | 20        | 1.0                      | 20.00     | 0           | 101                                   | 70       | 130         |      |          |      |
| Trichloroethene (TCE)       | 18        | 1.0                      | 20.00     | 0           | 90.3                                  | 70       | 130         |      |          |      |
| Surr: 1,2-Dichloroethane-d4 | 10        |                          | 10.00     |             | 99.6                                  | 70       | 130         |      |          |      |
| Surr: 4-Bromofluorobenzene  | 11        |                          | 10.00     |             | 112                                   | 70       | 130         |      |          |      |
| Surr: Dibromofluoromethane  | 10        |                          | 10.00     |             | 100                                   | 70       | 130         |      |          |      |
| Surr: Toluene-d8            | 11        |                          | 10.00     |             | 106                                   | 70       | 130         |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
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RL Reporting Detection Limit  
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30-Jun-16

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|                             |                  |     |                                 |             |  |                    |           |      |          |      |
|-----------------------------|------------------|-----|---------------------------------|-------------|--|--------------------|-----------|------|----------|------|
| Sample ID                   | <b>mb-25437</b>  |     | SampType: <b>MBLK</b>           |             | TestCode: <b>EPA Method 8270C: Semivolatiles</b> |                    |           |      |          |      |
| Client ID:                  | <b>PBW</b>       |     | Batch ID: <b>25437</b>          |             | RunNo: <b>34450</b>                              |                    |           |      |          |      |
| Prep Date:                  | <b>5/23/2016</b> |     | Analysis Date: <b>5/24/2016</b> |             | SeqNo: <b>1062341</b>                            | Units: <b>µg/L</b> |           |      |          |      |
| Analyte                     | Result           | PQL | SPK value                       | SPK Ref Val | %REC   | LowLimit           | HighLimit | %RPD | RPDLimit | Qual |
| Acenaphthene                | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Acenaphthylene              | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Aniline                     | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Anthracene                  | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Azobenzene                  | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Benz(a)anthracene           | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Benzo(a)pyrene              | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Benzo(b)fluoranthene        | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Benzo(g,h,i)perylene        | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Benzo(k)fluoranthene        | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Benzoic acid                | 6.7              | 20  |                                 |             |  |                    |           |      |          | J    |
| Benzyl alcohol              | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Bis(2-chloroethoxy)methane  | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Bis(2-chloroethyl)ether     | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Bis(2-chloroisopropyl)ether | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Bis(2-ethylhexyl)phthalate  | 3.8              | 10  |                                 |             |  |                    |           |      |          | J    |
| 4-Bromophenyl phenyl ether  | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Butyl benzyl phthalate      | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Carbazole                   | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 4-Chloro-3-methylphenol     | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 4-Chloroaniline             | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2-Chloronaphthalene         | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2-Chlorophenol              | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 4-Chlorophenyl phenyl ether | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Chrysene                    | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Di-n-butyl phthalate        | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Di-n-octyl phthalate        | 4.8              | 10  |                                 |             |  |                    |           |      |          | J    |
| Dibenz(a,h)anthracene       | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Dibenzofuran                | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 1,2-Dichlorobenzene         | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 1,3-Dichlorobenzene         | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 1,4-Dichlorobenzene         | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 3,3'-Dichlorobenzidine      | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Diethyl phthalate           | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Dimethyl phthalate          | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2,4-Dichlorophenol          | ND               | 20  |                                 |             |  |                    |           |      |          |      |
| 2,4-Dimethylphenol          | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 4,6-Dinitro-2-methylphenol  | ND               | 20  |                                 |             |  |                    |           |      |          |      |
| 2,4-Dinitrophenol           | ND               | 20  |                                 |             |  |                    |           |      |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                            |                  |     |                                 |             |  |                    |           |      |          |      |
|----------------------------|------------------|-----|---------------------------------|-------------|--|--------------------|-----------|------|----------|------|
| Sample ID                  | <b>mb-25437</b>  |     | SampType: <b>MBLK</b>           |             | TestCode: <b>EPA Method 8270C: Semivolatiles</b> |                    |           |      |          |      |
| Client ID:                 | <b>PBW</b>       |     | Batch ID: <b>25437</b>          |             | RunNo: <b>34450</b>                              |                    |           |      |          |      |
| Prep Date:                 | <b>5/23/2016</b> |     | Analysis Date: <b>5/24/2016</b> |             | SeqNo: <b>1062341</b>                            | Units: <b>µg/L</b> |           |      |          |      |
| Analyte                    | Result           | PQL | SPK value                       | SPK Ref Val | %REC   | LowLimit           | HighLimit | %RPD | RPDLimit | Qual |
| 2,4-Dinitrotoluene         | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2,6-Dinitrotoluene         | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Fluoranthene               | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Fluorene                   | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Hexachlorobenzene          | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Hexachlorobutadiene        | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Hexachlorocyclopentadiene  | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Hexachloroethane           | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Indeno(1,2,3-cd)pyrene     | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Isophorone                 | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 1-Methylnaphthalene        | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2-Methylnaphthalene        | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2-Methylphenol             | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 3+4-Methylphenol           | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| N-Nitrosodi-n-propylamine  | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| N-Nitrosodimethylamine     | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| N-Nitrosodiphenylamine     | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Naphthalene                | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2-Nitroaniline             | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 3-Nitroaniline             | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 4-Nitroaniline             | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Nitrobenzene               | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2-Nitrophenol              | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 4-Nitrophenol              | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Pentachlorophenol          | ND               | 20  |                                 |             |  |                    |           |      |          |      |
| Phenanthrene               | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Phenol                     | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Pyrene                     | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Pyridine                   | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 1,2,4-Trichlorobenzene     | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2,4,5-Trichlorophenol      | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| 2,4,6-Trichlorophenol      | ND               | 10  |                                 |             |  |                    |           |      |          |      |
| Surr: 2-Fluorophenol       | 120              |     | 200.0                           |             | 57.6   | 15                 | 123       |      |          |      |
| Surr: Phenol-d5            | 87               |     | 200.0                           |             | 43.6   | 4.13               | 124       |      |          |      |
| Surr: 2,4,6-Tribromophenol | 160              |     | 200.0                           |             | 79.7   | 18.4               | 134       |      |          |      |
| Surr: Nitrobenzene-d5      | 75               |     | 100.0                           |             | 74.9   | 28.8               | 134       |      |          |      |
| Surr: 2-Fluorobiphenyl     | 71               |     | 100.0                           |             | 71.3   | 35.9               | 125       |      |          |      |
| Surr: 4-Terphenyl-d14      | 65               |     | 100.0                           |             | 65.5   | 15                 | 146       |      |          |      |

### Qualifiers:

|    |   |    |   |
|----|---|----|---|
| *  | Value exceeds Maximum Contaminant Level.              | B  | Analyte detected in the associated Method Blank           |
| D  | Sample Diluted Due to Matrix                          | E  | Value above quantitation range                            |
| H  | Holding times for preparation or analysis exceeded    | J  | Analyte detected below quantitation limits                |
| ND | Not Detected at the Reporting Limit                   | P  | Sample pH Not In Range                                    |
| R  | RPD outside accepted recovery limits                  | RL | Reporting Detection Limit                                 |
| S  | % Recovery outside of range due to dilution or matrix | W  | Sample container temperature is out of limit as specified |

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

Client: Western Refining Company

Project: SWMU 10

|                            |           |     |                          |             |      |   |           |             |          |      |
|----------------------------|-----------|-----|--------------------------|-------------|------|---|-----------|-------------|----------|------|
| Sample ID                  | Ics-25437 |     | SampType: LCS            |             |      | TestCode: EPA Method 8270C: Semivolatiles |           |             |          |      |
| Client ID:                 | LCSW      |     | Batch ID: 25437          |             |      | RunNo: 34450                              |           |             |          |      |
| Prep Date:                 | 5/23/2016 |     | Analysis Date: 5/24/2016 |             |      | SeqNo: 1062342                            |           | Units: µg/L |          |      |
| Analyte                    | Result    | PQL | SPK value                | SPK Ref Val | %REC | LowLimit                                  | HighLimit | %RPD        | RPDLimit | Qual |
| Acenaphthene               | 84        | 10  | 100.0                    | 0           | 83.7 | 35  | 113       |             |          |      |
| 4-Chloro-3-methylphenol    | 130       | 10  | 200.0                    | 0           | 66.0 | 40.7                                      | 114       |             |          |      |
| 2-Chlorophenol             | 130       | 10  | 200.0                    | 0           | 64.4 | 37.6                                      | 113       |             |          |      |
| 1,4-Dichlorobenzene        | 69        | 10  | 100.0                    | 0           | 68.7 | 37.7                                      | 106       |             |          |      |
| 2,4-Dinitrotoluene         | 70        | 10  | 100.0                    | 0           | 70.1 | 37  | 91        |             |          |      |
| N-Nitrosodi-n-propylamine  | 74        | 10  | 100.0                    | 0           | 74.5 | 45.4                                      | 105       |             |          |      |
| 4-Nitrophenol              | 73        | 10  | 200.0                    | 0           | 36.4 | 33.4                                      | 104       |             |          |      |
| Pentachlorophenol          | 120       | 20  | 200.0                    | 0           | 58.0 | 29.5                                      | 94.9      |             |          |      |
| Phenol                     | 77        | 10  | 200.0                    | 0           | 38.4 | 30.6                                      | 119       |             |          |      |
| Pyrene                     | 82        | 10  | 100.0                    | 0           | 82.0 | 26.2                                      | 120       |             |          |      |
| 1,2,4-Trichlorobenzene     | 82        | 10  | 100.0                    | 0           | 81.6 | 39.9                                      | 125       |             |          |      |
| Surr: 2-Fluorophenol       | 95        |     | 200.0                    |             | 47.6 | 15  | 123       |             |          |      |
| Surr: Phenol-d5            | 75        |     | 200.0                    |             | 37.5 | 4.13                                      | 124       |             |          |      |
| Surr: 2,4,6-Tribromophenol | 150       |     | 200.0                    |             | 74.3 | 18.4                                      | 134       |             |          |      |
| Surr: Nitrobenzene-d5      | 68        |     | 100.0                    |             | 68.2 | 28.8                                      | 134       |             |          |      |
| Surr: 2-Fluorobiphenyl     | 70        |     | 100.0                    |             | 70.0 | 35.9                                      | 125       |             |          |      |
| Surr: 4-Terphenyl-d14      | 61        |     | 100.0                    |             | 61.5 | 15  | 146       |             |          |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified

# QC SUMMARY REPORT

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1605998

30-Jun-16

**Client:** Western Refining Company

**Project:** SWMU 10

| Sample ID <b>rb1</b>          | SampType: <b>MBLK</b>           |       |           | TestCode: <b>EPA Method 8015D: Gasoline Range</b> |      |          |                    |      |          |      |
|-------------------------------|---------------------------------|-------|-----------|---|------|----------|--------------------|------|----------|------|
| Client ID: <b>PBW</b>         | Batch ID: <b>C34513</b>         |       |           | RunNo: <b>34513</b>                               |      |          |                    |      |          |      |
| Prep Date:                    | Analysis Date: <b>5/26/2016</b> |       |           | SeqNo: <b>1064532</b>                             |      |          | Units: <b>mg/L</b> |      |          |      |
| Analyte                       | Result                          | PQL   | SPK value | SPK Ref Val                                       | %REC | LowLimit | HighLimit          | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO) | ND                              | 0.050 |           |   |      |          |                    |      |          |      |
| Surr: BFB                     | 9.7                             |       | 10.00     |   | 96.9 | 70       | 130                |      |          |      |

| Sample ID <b>2.5ug gro lcs</b> | SampType: <b>LCS</b>            |       |           | TestCode: <b>EPA Method 8015D: Gasoline Range</b> |      |          |                    |      |          |      |
|--------------------------------|---------------------------------|-------|-----------|---|------|----------|--------------------|------|----------|------|
| Client ID: <b>LCSW</b>         | Batch ID: <b>C34513</b>         |       |           | RunNo: <b>34513</b>                               |      |          |                    |      |          |      |
| Prep Date:                     | Analysis Date: <b>5/26/2016</b> |       |           | SeqNo: <b>1064533</b>                             |      |          | Units: <b>mg/L</b> |      |          |      |
| Analyte                        | Result                          | PQL   | SPK value | SPK Ref Val                                       | %REC | LowLimit | HighLimit          | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO)  | 0.52                            | 0.050 | 0.5000    | 0   | 103  | 75.4     | 118                |      |          |      |
| Surr: BFB                      | 10                              |       | 10.00     |   | 101  | 70       | 130                |      |          |      |

| Sample ID <b>1605998-001B MS</b> | SampType: <b>MS</b>             |      |           | TestCode: <b>EPA Method 8015D: Gasoline Range</b> |      |          |                    |      |          |      |
|----------------------------------|---------------------------------|------|-----------|---|------|----------|--------------------|------|----------|------|
| Client ID: <b>SWMU 10-20-GW</b>  | Batch ID: <b>C34513</b>         |      |           | RunNo: <b>34513</b>                               |      |          |                    |      |          |      |
| Prep Date:                       | Analysis Date: <b>5/26/2016</b> |      |           | SeqNo: <b>1064538</b>                             |      |          | Units: <b>mg/L</b> |      |          |      |
| Analyte                          | Result                          | PQL  | SPK value | SPK Ref Val                                       | %REC | LowLimit | HighLimit          | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO)    | 29                              | 0.50 | 5.000     | 25.03   | 72.5 | 53.8     | 128                |      |          |      |
| Surr: BFB                        | 97                              |      | 100.0     |   | 97.1 | 70       | 130                |      |          |      |

| Sample ID <b>1605998-001B MSD</b> | SampType: <b>MSD</b>            |      |           | TestCode: <b>EPA Method 8015D: Gasoline Range</b> |      |          |                    |      |          |      |
|-----------------------------------|---------------------------------|------|-----------|---|------|----------|--------------------|------|----------|------|
| Client ID: <b>SWMU 10-20-GW</b>   | Batch ID: <b>C34513</b>         |      |           | RunNo: <b>34513</b>                               |      |          |                    |      |          |      |
| Prep Date:                        | Analysis Date: <b>5/26/2016</b> |      |           | SeqNo: <b>1064539</b>                             |      |          | Units: <b>mg/L</b> |      |          |      |
| Analyte                           | Result                          | PQL  | SPK value | SPK Ref Val                                       | %REC | LowLimit | HighLimit          | %RPD | RPDLimit | Qual |
| Gasoline Range Organics (GRO)     | 29                              | 0.50 | 5.000     | 25.03   | 78.6 | 53.8     | 128                | 1.05 | 20       |      |
| Surr: BFB                         | 98                              |      | 100.0     |   | 98.4 | 70       | 130                | 0    | 0        |      |

### Qualifiers:

\* Value exceeds Maximum Contaminant Level.  
D Sample Diluted Due to Matrix  
H Holding times for preparation or analysis exceeded  
ND Not Detected at the Reporting Limit  
R RPD outside accepted recovery limits  
S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank  
E Value above quantitation range  
J Analyte detected below quantitation limits  
P Sample pH Not In Range  
RL Reporting Detection Limit  
W Sample container temperature is out of limit as specified



Hall Environmental Analysis Laboratory  
4901 Hawks NE  
Albuquerque, NM 87109  
TEL: 505-345-3975 FAX: 505-345-4107  
Website: www.hallenvironmental.com

## Sample Log-In Check List

Client Name: Western Refining Gallup

Work Order Number: 1605998

RcptNo: 1

Received by/date: *AF* *05/20/16*

Logged By: Lindsay Mangin 5/20/2016 3:00:00 PM *Judy Mangin*

Completed By: Lindsay Mangin 5/20/2016 3:34:27 PM *Judy Mangin*

Reviewed By: *JA* *05/23/16*

### Chain of Custody

1. Custody seals intact on sample bottles? Yes ☒ No ☐ Not Present ☐  
2. Is Chain of Custody complete? Yes ☒ No ☐ Not Present ☐  
3. How was the sample delivered? Courier

### Log In

4. Was an attempt made to cool the samples? Yes ☒ No ☐ NA ☐  
5. Were all samples received at a temperature of  $>0^{\circ}\text{C}$  to  $6.0^{\circ}\text{C}$ ? Yes ☒ No ☐ NA ☐  
6. Sample(s) in proper container(s)? Yes ☒ No ☐  
7. Sufficient sample volume for indicated test(s)? Yes ☒ No ☐  
8. Are samples (except VOA and ONG) properly preserved? Yes ☒ No ☒  
9. Was preservative added to bottles? Yes ☒ No ☒ NA ☐  
10. VOA vials have zero headspace? Yes ☒ No ☐ No VOA Vials ☐  
11. Were any sample containers received broken? Yes ☐ No ☒  
12. Does paperwork match bottle labels?  
(Note discrepancies on chain of custody) Yes ☒ No ☐  
13. Are matrices correctly identified on Chain of Custody? Yes ☒ No ☐  
14. Is it clear what analyses were requested? Yes ☒ No ☐  
15. Were all holding times able to be met?  
(If no, notify customer for authorization.) Yes ☒ No ☐

# of preserved bottles checked for pH: *2,1*  
Adjusted? *yes*  
Checked by: *AS*

### Special Handling (if applicable)

16. Was client notified of all discrepancies with this order? Yes ☐ No ☐ NA ☒

Person Notified: \_\_\_\_\_ Date: \_\_\_\_\_  
By Whom: \_\_\_\_\_ Via: ☐ eMail ☐ Phone ☐ Fax ☐ In Person  
Regarding: \_\_\_\_\_  
Client Instructions: \_\_\_\_\_

17. Additional remarks: *For Metals Analysis. Added 1 mL HNO<sub>3</sub> to -001F for acceptable pH, held for 24hrs prior to analysis. AG 05/23/16*  
18. Cooler Information

| Cooler No | Temp °C | Condition | Seal Intact | Seal No | Seal Date | Signed By |
|-----------|---------|-----------|-------------|---------|-----------|-----------|
| 1         | 2.2     | Good      | Yes         |         |           |           |

# Chain-of-Custody Record

Client: WESTERN REFINING SW, INC.

GALLUP REFINERY

Mailing Address: 922 GIANT CROSSING RD

GALLUP, NM 87301

Phone #: 505-722-0231

email or Fax#: ED.RIESE@WNR.COM

QA/QC Package:

☐ Standard

☒ Level 4 (Full Validation)

Accreditation

☐ NELAP

☐ Other

☒ EDD (Type)

EXCEL

Sample Temperature: 2.2

Container Type and #

Preservative Type

HEAL No.

Date

Time

Matrix

Sample Request ID

5/20/16 1045 WATER SWMU 10-20-GW

1 LITER

AMBER-2

250 ML

AMBER-1

100 ML

PLASTIC-1

500 ML

PLASTIC-1

50 ML

PLASTIC-1

500 ML

PLASTIC-1

40 ML

VOA-4

HCL

NEAT

500 ML

PLASTIC-1

NEAT

500 ML

PLASTIC-1

40 ML

VOA-6

HCL

NEAT

1 LITER

AMBER-1

500 ML

PLASTIC-1

NEAT

500 ML

PLASTIC-1

40 ML

VOA-4

HCL

NEAT

500 ML

PLASTIC-1

NEAT

500 ML

PLASTIC-1

40 ML

VOA-4

HCL

NEAT

500 ML

PLASTIC-1

NEAT

500 ML

PLASTIC-1

40 ML

VOA-4

HCL

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**WESTERN REFINING SOUTHWEST, INC.**  
**GALLUP REFINERY - NORTH DRAINAGE DITCH**  
**METALS AND CYANIDE ANALYSES FOR GROUNDWATER SAMPLES**  
**AND WATER QA/QC SAMPLES**  
**TOTAL METALS ANALYSIS AND DISSOLVED METALS ANALYSIS**

| <b>Analyte</b> | <b>Analytical Method</b>      |
|----------------|-------------------------------|
| Antimony       | SW-846 method 6010/6020       |
| Arsenic        | SW-846 method 6010/6020       |
| Barium         | SW-846 method 6010/6020       |
| Beryllium      | SW-846 method 6010/6020       |
| Cadmium        | SW-846 method 6010/6020       |
| Chromium       | SW-846 method 6010/6020       |
| Cobalt         | SW-846 method 6010/6020       |
| Cyanide        | SW-846 method 335.4/335.2 mod |
| Lead           | SW-846 method 6010/6020       |
| Mercury        | SW-846 method 7470/7471       |
| Nickel         | SW-846 method 6010/6020       |
| Selenium       | SW-846 method 6010/6020       |
| Silver         | SW-846 method 6010/6020       |
| Vanadium       | SW-846 method 6010/6020       |
| Zinc           | SW-846 method 6010/6020       |
| Iron           | SW-846 method 6010/6020       |
| Manganese      | SW-846 method 6010/6020       |

**GENERAL CHEMISTRY PARAMETERS FOR GROUNDWATER SAMPLES**  
**AND WATER QA/QC SAMPLES**

| <b>Analyte</b> | <b>Analytical Method</b> |
|----------------|--------------------------|
| Chloride       | EPA method 300.0         |
| Fluoride       | EPA method 300.0         |
| Sulfate        | EPA method 300.0         |

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## **Appendix E**

### **Quality Assurance/Quality Control Review**

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## DATA VALIDATION INTRODUCTION

This summary presents data verification results for soil and groundwater samples collected from soil boring and monitoring wells installed at SWMU10 at the Gallup Refinery. The data review was performed in accordance with Provision IV.J.3.b (Review of Field and Laboratory QA/QC Data) of the RCRA Permit issued by NMED in October 2013, USEPA Functional Guidelines for Organic and Inorganic Data Review, and quality assurance and control parameters set by the project laboratory Hall Environmental Analysis Laboratory, Inc.

A total of 73 soil samples and 11 groundwater samples (excluding QA samples) were collected from April 28, 2015 through September 21, 2016 in accordance with the SWMU 10 Investigation Work Plan (Western Refining Southwest, Inc., 2014). Soil and groundwater samples were submitted to Hall Environmental Analysis Laboratory for the following parameters in accordance with the approved Work Plan:

- volatile organic compounds (VOCs) by USEPA Method 8260B;
- semi-volatile organic compounds (SVOCs) by USEPA Method 8270;
- Gasoline, diesel, and motor oil range organics by SW-846 Method 8015B;
- Total recoverable and dissolved metals (Antimony, arsenic, barium, beryllium, cadmium, chromium, cobalt, lead, nickel, selenium, silver, vanadium, and zinc) by SW846 Method 6010/6020;
- Chromium VI by SW-846 method 3060A;
- Tetraethyl lead by SW-846 method 3546;
- Cyanide by SW-846 method 9012; and
- Mercury by EPA Method 7470.

The groundwater samples were analyzed for water quality parameters including, sulfate, chloride, and fluoride.

Additionally, 29 quality assurance samples consisting of trip blanks, field blanks, equipment rinsate blanks, and field duplicates were collected and analyzed as part of the investigation activities. Table A-1 presents a summary of the field sample identifications, laboratory sample identifications, and sample collection dates.

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## QUALITY CONTROL PARAMETERS REVIEWED

Sample results were subject to a Level II data review that includes an evaluation of the following quality control (QC) parameters:

- Chain-of-Custody;
- Sample Preservation and Temperature Upon Laboratory Receipt
- Holding Times;
- Blank Contamination (method blanks, trip blanks, field blanks, and equipment rinsate blanks);
- Surrogate Recovery (for organic parameters);
- Laboratory Control Sample (LCS) Recovery and Relative Percent Difference (RPD);
- Matrix Spike/Matrix Spike Duplicate (MS/MSD) Recovery and RPD;
- Duplicates (field duplicate, laboratory duplicate); and
- Other Applicable QC Parameters.

The data qualifiers used to qualify the analytical results associated with QC parameters outside of the established data quality objectives are defined below:

- |    |  |
|----|--|
| J+ | The analyte was positively identified; however, the result should be considered an estimated value with a potential high bias. |
| J- | The analyte was positively identified; however, the result should be considered an estimated value with a potential low bias.  |
| UJ | The reporting limit for a constituent that was not detected is considered an estimated value.                                  |
| R  | Quality control indicates that the data is not usable.   |

Results qualified as “J+”, “J-”, or “UJ” are of acceptable data quality and may be used quantitatively to fulfill the objectives of the analytical program, per EPA guidelines.

Results for the performance monitoring events that required qualification based on the data verification are summarized in Table A-2.

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## **CHAIN-OF-CUSTODY**

The chain-of-custody documentation associated with project samples was found to be complete. Chain-of-custodies included sample identifications, date and time of collection, requested parameters, and relinquished/received signatures.

## **SAMPLE PRESERVATION AND TEMPERATURE UPON LABORATORY RECEIPT**

Samples collected were received preserved and intact by Hall Environmental Laboratories, Inc. Samples were received by the laboratory at a temperature of 6.0 degrees Celsius or lower. Data qualification on lower temperature samples was not required. The hexavalent chromium field sample SWMU 10-12-GW was received at a pH of 7.5 rather than the range of 8 to 10 units. The field sample was not qualified. The sample bottle for the total metals analyses for sample SWMU 10-20-GW was apparently received with a pH outside the desired range, thus the lab added one milliliter of HNO<sub>3</sub> and held the sample for 24 hours prior to analysis to ensure no metals had precipitated in the sample. The associated sample results are not qualified.

## **HOLDING TIMES**

All samples were extracted and analyzed within method-specified holding time limits with the exception of the following:

- Lab Report 1505618 - Sample SWMU 10 EB06 was extracted 2 days after the 7 day holding time. This affected only the Method 8270C semi volatile analyses and only one sample. Since results may be biased low, the sample non-detect results were qualified "UJ".

See Table A-2 for qualified data.

## **BLANK CONTAMINATION**

### **Method Blank**

Method blanks were analyzed at the appropriate frequency. Target compounds were not detected in the method blanks, with the exception of the following:

- 1,2,4-Trimethylbenzene, acetone, 2-butanone (MEK), chloromethane, and 1,1-dichloroethane were detected in the blank for batch 19008. Acetone, 2-butanone, chloromethane, and 1,1-dichloroethane were not detected in the associated samples and qualification was not required. 1,2,4-Trimethylbenzene was detected at concentrations
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greater than 5x blank in associated samples; and qualification of the field samples was not necessary.

- Zinc was detected in the method blank for batch 19025. Zinc was detected at similar concentrations in associated sample SWMU 10-3-GW; and the field sample was qualified with a "J+".
  - Toluene and 2-butanone were detected in the blank for batch 19032. Toluene was detected in 2 associated samples and 2-butanone was detected in 10 associated samples at similar concentrations; and the field samples were qualified with a "J+".
  - Lead, iron and zinc were detected in the method blank for batch for 19080. Lead, iron and zinc were detected in associated samples at concentrations greater than 5x the concentrations in the blank; and qualification of field samples was not necessary;
  - Lead and iron were detected in the method blank for batch 19081. Lead and iron were detected in associated samples. The concentration for lead in one sample SWMU 10-8 (18-20') was less than 5x the concentration in the blank and this sample was qualified "J+". The results for remaining associated samples were greater than 5x the blank and qualification was not necessary.
  - Cadmium and iron were detected in the method blank for batch 19082. Cadmium was not detected in associated samples. Iron was detected in associated samples but the concentrations were greater than 5x the blank and qualification of field samples was not necessary.
  - 2-Butanone and toluene were detected in the method blank for batch 19107. 2-Butanone was detected in six associated samples at similar concentrations; and the field samples were qualified "J+". Toluene was detected in associated samples at concentrations greater than 5x the blank and qualification of the field samples was not necessary.
  - 2-Butanone was detected in the blank for batch 19243. 2-Butanone was detected in three associated samples SWMU 10-16 (2-4'), SWMU 10-16 (4.5-5') and SWMU 10-16 (8-9") at similar concentrations; and the field samples were qualified "J+".
  - Iron and silver were detected in the method blank for batch 19259. Silver was not detected in associated samples. Iron was detected in associated samples but the concentrations were greater than 5x the blank; and field samples were not qualified.
  - Iron was detected in the blank for batch 19279. Iron was detected in associated samples but the concentrations were greater than 5x the blank and field samples were not qualified.
  - Iron was detected in the blank for batch 19280. Iron was detected in associated samples but the concentrations were greater than 5x the blank and field samples were not qualified.
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- Iron was detected in the blank for batch 19333. Iron was detected in associated samples but the concentrations were greater than 5x the blank and field samples were not qualified.
  - Mercury was detected in the blank for batch 19401. Mercury was also detected in the equipment blank EB06 and in an associated sample SWMU 10-11-GW at similar concentrations. The field samples were qualified "J+".
  - Beryllium was detected in the method blank for batch R26030. Beryllium was detected in three associated samples SWMU 10-1-GW, SWMU 10-3-GW and SWMU 10-5-GW at similar concentrations. The field samples were qualified "J+".
  - 2-Butanone, chloroethane, and chloromethane were detected in the method blank for batch R26144. 2-Butanone, chloroethane, and chloromethane were not detected in associated samples and qualification of field samples is not necessary.
  - 1,2,4- Trimethylbenzene, 2-butanone, and toluene were detected in the method blank for batch R25982. 1,2,4- Trimethylbenzene was detected in associated samples SWMU 10-1 (2-4') and SWMU 10-8 (4-6') at a similar concentration and the field samples were flagged "J+". 2-Butanone was not detected and toluene was detected in associated samples but concentrations were 5x the blank; and qualification of field samples is not necessary.
  - 1,2,4-Trimethylbenzene, 1,2-dichloroethane, bromomethane, and chloroethane were detected in the blank for batch R26221. 1,2,4-Trimethylbenzene was detected in associated samples but concentrations were 5x the blank; and bromomethane, chloroethane, and 1,2-Dichloroethane were not detected in associated samples. Qualification of field samples was not necessary.
  - Silver was detected in the method blank for analytical batch R25881. Silver was not detected in the associated sample and qualification was not required.
  - Bromomethane and 1,2,4-trimethylbenzene were detected in the method blank for batch R26209; Bromomethane was detected in associated samples SWMU 10-11 (8-10"), SWMU 10-13 (6-8') and SWMU 10-14 (6-8') at similar concentrations; and field samples were flagged "J+". 1,2,4-Trimethylbenzene was detected in associated samples but concentrations were 5x the blank; and qualification of field samples was not necessary.
  - Barium was detected in the blank for batch R26242. Barium was detected in associated samples but concentrations were 5x the blank; and qualification of field samples is not necessary.
  - Silver was detected in the blank for batch R26291. Silver was not detected in associated samples; and qualification of field samples was not necessary.
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- 1,2,4-Trimethylbenzene, 1,2-dichlorethane, and vinyl chloride were detected in the blank for batch R26332. 1,2,4-Trimethylbenzene was detected in associated samples SWMU 10-12-GW, SWMU 10-14-GW, SWMU 10-16-GW and two Trip Blanks at similar concentrations; and field samples were flagged “J+”. 1,2-Dichlorethane and vinyl chloride were detected in associated sample SWMU 10-16-GW at similar concentrations and field samples were flagged “J+”.
  - Bis(2-ethylhexyl)phthalate, di-n-butyl phthalate, and 4,6-dinitro-2-methylphenol were detected in the method blank for batch 25469 but many sample results were non-detect and are not qualified. Samples SWMU 10-18 (2-2.5' and 8-10') had detections of di-n-butyl phthalate at concentrations close to that detected in the method blank and are both flagged as “J+”. Samples SWMU 10-21 (2-2.5', 12-14', and 20-22') and SWMU 10-22 (2-2.5' and 8-9') were flagged as J+ for bis(2-ethylhexyl)phthalate. Samples SWMU 10-21 (12-14') and SWMU 10-22 (2-2.5' and 8-9') were flagged as J+ for di-n-butyl phthalate.
  - Mercury was detected in the method blank for sample batch 25448 and associated samples SWMU 10-18 (2-2.5', 8-10', and 18-20'), SWMU 10-19 (2-2.5', 12-14' and 18-20'), and SWMU 10-20 (2-2.5', 8-10', 10-12', 16-18', and 20-22') are all qualified as J+.
  - Manganese was detected in the method blank for sample batch 25447 and associated samples SWMU 10-18 (2-2.5', 8-10', and 18-20') are all qualified as J+.
  - 1-Methlynaphthalene and methylene chloride were both detected in the method blank for batch D34413. Associated samples SWMU 10-19 (12-14' and 18-20') and SWMU 10-20 (8-10', 10-12', and 16-18') all flagged as J+ for 1-methlynaphthalene, while only samples SWMU 10-19 (12-14') and SWMU 10-20 (8-10' and 10-12') are flagged as J+ for methylene chloride.
  - Cyanide, mercury and methylene chloride were detected in the method blanks for batches WG876094, 25512, and A34513, respectively and the associated results in sample SWMU 10-20-GW are flagged as J+.
  - Toluene and bromomethane were detected in the method blank for batch S37464 and associated sample SWMU 10-24 (6-8') is flagged as J+.
  - 1,2,4-Trimethylbenzene was detected in the method blank for batch S37484 and sample SWMU 10-25 (2-2.5') is flagged as J+.
  - 1-Methlynaphthalene was detected in method blank for batch S37484 and samples SWMU 10-23 (15-16'), SWMU 10-24 (15-16'), and SWMU 10-25 (2-2.5') were flagged as J+.
  - Bis(2-ethylhexyl)phthalate and di-n-butyl phthalate were both detected in the method blank for batch W37464 and associated samples SWMU 10-23 (2-2.5' and 15-16'), SWMU 10-24
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(2-2.5', 8-10', and 15-16'), and SWMU 10-25 (2-2.5', 10-12', and 16.8-18') are flagged as J+.

- Manganese and zinc were detected in the method blank for batch 27709 and associated samples SWMU 10-23 (2-2.5' and 15-16'), SWMU 10-24 (2-2.5', 6-8', 8-10', and 15-16'), and SWMU 10-25 (2-2.5', 10-12', and 16.8-18') are flagged as J+.
- Manganese was detected in the method blank for batch 27858 and associated samples SWMU 10-25-GW is flagged as J+. In addition, bis(2-ethylhexyl)phthalate was detected in the batch (27668) that is associated with SWMU 10-25-GW and the result is flagged as J+.

See Table A-2 for qualified data.

### **Trip Blank**

Trip blanks were analyzed at the appropriate frequency as specified in the Permit. Target compounds were not detected in the trip blanks with the following exceptions:

#### **Lab Report 1505698**

- The VOCs ethylbenzene (0.11-J ug/L), naphthalene (0.31-J ug/L), 1-methylnaphthalene (0.71-J ug/L), 2-methylnaphthalene (1.1-J ug/L), 1,2,4-trimethylbenzene (0.29-J ug/L), and xylenes (0.041-J ug/L) were detected in TRIP Blank Sample No. 1501698-002a. The analytes were detected in the groundwater sample 1505698-001a associated with this trip blank at concentrations more than 5x the concentration detected in the blank. Sample results were not qualified; and

#### **Lab Report 1505700**

- The VOCs toluene (0.14-J ug/L) and 1,2,4-trimethylbenzene (0.18-J ug/L) were detected in TRIP Blank Sample No. 1505700-002a. Toluene was not detected and 1,2,4-trimethylbenzene was detected in the groundwater sample 1505700-001a (SWMU 10-12-GW) associated with this trip blank. The detected concentration of 1,2,4-trimethylbenzene (0.66 ug/L) was less than 5 times the concentration in the trip blank and may be biased high. The 1,2,4-trimethylbenzene result for the field sample was flagged "J+".

See Table A-2 for qualified data.

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### **Field Blanks/Equipment Rinsate Blank**

Field and equipment rinsate blanks were collected as specified in the SWMU10 Investigation Work Plan and the Permit.

### **Common Laboratory Contaminants**

Per USEPA guidelines, common laboratory contaminants for VOC analysis are acetone, 2-butanone (MEK), cyclohexane, chloromethane, and methylene chloride. Common laboratory contaminants for SVOC analysis include phthalates. Data qualification was not required for MEK in sample batches 19032, 19107, and 19243 where the laboratory contaminant was detected in the blank. Data qualification was required for other samples with methylene chloride and phthalates since there were detection in blanks and field analytical results were detected at concentrations less than 10 times the blank concentration in field samples. See Table A-2 for qualified data.

### **Methanol Blanks**

Methanol Blanks provided by the laboratory were analyzed for VOCs. There were no analytes detected in the methanol blanks above the respective laboratory reporting limits.

### **SURROGATE RECOVERY**

Surrogate recoveries for the organic and inorganic analyses were performed at the required frequency and were within laboratory acceptance limits, with the following exceptions:

Lab Report 1504C86

- Surrogate recovery for nitrobenzene-d6 was high and above the acceptance limit for field sample SWMU 10 EB01. The surrogate recovery for five of the six surrogates included in Method 8270C were within limits. Since only one surrogate was outside limits, the associated field sample results for semi-volatile organic compounds were not qualified.

Lab Report 1505002

- Surrogate recovery for nitrobenzene-d6 was high and above the acceptance limit for field sample SWMU 10 EB02. The surrogate recovery for five of the six surrogates included in Method 8270C were within limits. Since only one surrogate was outside limits, the associated field sample results for semi-volatile organic compounds were not qualified.
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#### Lab Report 1505003

- Surrogate recovery for di-n-octyl phthalate (DNOP) was below the acceptance limit for field samples SWMU 10-5 (0-2') and SWMU 10-5 (4-6'). Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8015 diesel range organics (DRO) or matrix effects; therefore data qualification was not required.
- Surrogate recoveries for Method 8270C were below the lower acceptance limits for field sample SWMU 10-5 (0-2'). Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8270C or matrix effects; therefore data qualification was not required.

#### Lab Report 1505004

- Surrogate recovery for DNOP was below the lower acceptance limit for field sample SWMU 10-4 (2-4'). Low surrogate was related to required sample dilution for analytical analysis by Method 8015 DRO or matrix effects; therefore data qualification was not required.
- Surrogate recoveries for Method 8270C were below the lower acceptance limits for field sample SWMU 10 DUP01. Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8270C or matrix effects; therefore data qualification was not required.

#### Lab Report 1505005

- Surrogate recovery for bromofluorobenzene (BFB) was high and above the upper acceptance limit for field sample SWMU 10-3-GW. The associated field sample results for Method 8015 gasoline range organics (GRO) are qualified J+ due to a potential high bias.

#### Lab Report 1505057

- Surrogate recovery for DNOP was below the lower acceptance limit for field sample SWMU 10-8 (2-4'). Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8015 DRO or matrix effects; therefore data qualification was not required.
  - Surrogate recoveries for Method 8270C were below the lower acceptance limits for field samples SWMU 10-9 (4-6') and SWMU 10-8 (2-4'). Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8270C or matrix effects; therefore data qualification was not required.
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#### Lab Report 1505058

- Surrogate recovery for DNOP was below the lower acceptance limit for field sample SWMU 10-10 (4-6'). Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8015 DRO or matrix effects; therefore data qualification was not required.
- Surrogate recoveries for Method 8270C were below the lower acceptance limits for field sample SWMU 10-10 (4-6'). Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8270C or matrix effects; therefore data qualification was not required.

#### Lab Report 1505218

- Surrogate recovery for BFB was high and above the upper acceptance limit for field sample SWMU 10-5-GW. The associated field sample results for Method 8015 GRO are qualified J+ due to a potential high bias. Non-detect results are not qualified.
- Surrogate recoveries for 2-fluorophenol and 2,4,6-tribromophenol were below the acceptance limit for field sample SWMU 10-5-GW. The surrogate recovery for three of the six surrogates included in Method 8270C were within limits. Since only two of six surrogates were outside limits, the associated field sample results for semi-volatile organic compounds were not qualified.

#### Lab Report 1505570

- Surrogate recovery for DNOP was below the lower acceptance limit for field samples SWMU 10-14 (6-8') and SWMU 10-11 (4-6'). Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8015 DRO or matrix interference; therefore data qualification was not qualified.
  - Surrogate recovery for DNOP was high and above the upper acceptance limit for field sample SWMU 10-11 (8-10'). The associated field sample results for Method 8015 GRO are qualified J+ due to a potential high bias. Non-detect results are not qualified.
  - Surrogate recoveries for Method 8270C were below the lower acceptance limits for field samples SWMU 10-9 (4-6') and SWMU 10-8 (2-4'). Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8270C or matrix effects; therefore data qualification was not required.
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#### Lab Report 15056174

- Surrogate recovery for DNOP was below the lower acceptance limit for field samples SWMU 10-13 (6-8'), SWMU 10-17 (6-8'), and SWMU 10 DUP04. Low surrogate recovery was due to required sample dilution for analytical analysis by Method 8015 DRO or matrix effects; therefore data qualification was not qualified.
- Surrogate recoveries for Method 8270C were below the lower acceptance limits for field samples SWMU 10-17 (6-8') and SWMU DUP04. Low surrogate recovery was related to required sample dilution for analytical analysis by Method 8270C or matrix effects; therefore data qualification was not required.

#### Lab Report 1505709

- Surrogate recovery for nitrobenzene d-5 was high and above the upper acceptance limit for field sample SWMU 10 DUP01GW. The surrogate recovery for five of the six surrogates included in Method 8270C were within limits. Since only one of six surrogates was outside limits, the associated field sample results for semi-volatile organic compounds were not qualified.
- Surrogate recovery for dibromofluorobenzene was high and above the acceptance limit for field sample SWMU10 DUP01GW. The surrogate recovery for four of the five surrogates included in Method 8260 were within limits. Since only one of five surrogates was outside limits, the associated field sample results for volatile organic compounds were not qualified.

See Table A-2 for qualified data.

#### **LCS RECOVERY AND RELATIVE PERCENT DIFFERENCE**

Laboratory control samples (LCS0/LCS duplicates) were performed at the required frequency and were evaluated based on the following criteria:

- If the analyte recovery was above acceptance limits for the LCS or LCS duplicate, but the analyte was not detected in the associated batch, then data qualification was not required.
  - If the analyte recovery was above acceptance limits for the LCS or LCS duplicate and the analyte was detected in the associated batch, then the analyte results were qualified "J+" to account for a potential high bias.
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- If the analyte recovery was below acceptance limits for LCS or LCS duplicate then the analyte results in the associated analytical batch were qualified (“UJ” for non-detects and “J-” for detected results) to account for a potential low bias.

LCS/LCSD percent recoveries and relative percent differences (RPDs) were within acceptance limits and no qualification was required.

#### **MS/MSD RECOVERY AND RELATIVE PERCENT DIFFERENCE**

Matrix Spike/Matrix Spike Duplicate (MS/MSD) samples were performed at the required frequency and were evaluated by the following criteria:

- If the MS or MSD recovery for an analyte was above acceptance limits but the analyte was not detected in the associated analytical batch, then data qualification was not required.
- If the MS or MSD recovery for an analyte was above acceptance limits and the analyte was detected in the associated analytical batch, then analyte results were qualified “J+” to account for a potential high bias.
- Low MS/MSD recoveries for organic or inorganic parameters result in sample qualification of the associated analytical batch with a “J-”.
- Results were not qualified based on non-project specific MS/MSD (i.e., batch QC) recoveries.

MS/MSD percent recoveries and RPDs were within acceptance limits except for the following:

##### **Lab Report 1505003**

- The MS/MSD recoveries for benzene and toluene for Method 8260B Batch 19032 were below the lower acceptance limits for field sample SWMU 10-5 (22-24). Low recovery was related to dilution or matrix interference. Since results may be biased low, detected results in the batch are qualified “J-” and non-detect results are qualified “UJ”.

##### **Lab Report 1505057**

- The MS recovery of 127% and the MSD recovery of 172% for mercury was above the acceptance limit of 125% for Method 7471 batch 19445 and field sample SWMU 10-9 (2-4'). Since results may be biased high, the detected results are qualified “J+”.
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#### Lab Report 1505218

- The MS/MSD recovery of 166% for BFB is above the upper limit of 120% for Method 8015 GRO batch R26016 in field sample SWMU 10-5-GW. The associated field sample results for Method 8015 GRO are qualified J+ due to a potential high bias. Non-detect samples were not qualified.
- The MS/MSD recoveries were low and the relative percent difference (RPD) were outside the acceptance limits for several chemicals and surrogates for Method 8270C batch 19150 performed for field sample SWMU 10-5-GW. Low recovery was related to dilution or matrix interference. Since results may be biased low, the associated semi-volatile results were qualified "J-"if detected or "UJ" if non-detect.

#### Lab Report 1505570

- The MS/MSD recoveries for all compounds and surrogates for Method 8270C batch 19270 were within limits for field sample SWMU 10-14 (4-6'). The RPD of 30.4 for N-nitroso-n-propyl amine was above the limit of 27.5. Since the RPDs for other compounds were within limits, and the N-nitroso-n-propyl amine RPD was only slightly out of specifications, the field data were not qualified.
- The MS recovery of 33.2% and the MSD recovery of 33.3% for antimony was below the lower limit of 75% for Method 6010 batch 19259 and field sample SWMU 10-14 (6-8'). Low recovery was related to dilution or matrix interference. The associated field sample results for antimony are qualified "UJ" due to a potential low bias.
- The MS recovery for lead was within limits and MSD recovery of 74.4 for lead was below the acceptance limit of 75% for Method 6010 batch 19259 and field sample SWMU 10-14 (6-8'). Since MSD recovery was only slightly low, the associated field sample results were not qualified.

#### Lab Report 1505698

- MS recoveries for benzene (16.4%), toluene (13.4%), chlorobenzene (11.7%), 1,1-dichloroethylene (6.58%), and trichloroethylene (9.98%) were below the lower acceptance limit of 30% for Method 8260 batch R26322 and field sample SWMU 10-11-GW. The RPDs for these compounds were outside acceptance limits. Low recovery was related to dilution or matrix interference. The field sample results for volatile organics are qualified "J-"and non-detect results are qualified "UJ" due to a potential low bias.
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#### Lab Report 1605874

- The MS and MSD recoveries for N-Nitrosodi-n-propylamine, 42.7% and 40.7, respectively, were below the lower limit of 43.5%. Antimony, cobalt, lead, manganese, nickel, and selenium had low MS recoveries of 29.8%, 69.2%, 67%, 74.6%, 72.6%, and 68.2%, respectively, vs the lower limit of 75%. Antimony, arsenic, beryllium, cadmium, cobalt, lead, manganese, nickel, and selenium had low MSD recoveries of 27.2%, 71.4%, 73.6%, 74.6%, 66.6%, 63.9%, 74.4%, 71.7%, and 66.2%, respectively, vs the lower limit of 75% .Low recovery was related to dilution or matrix interference. The field sample results with low recovery are qualified “J-“and non-detect results are qualified “UJ” due to a potential low bias. Barium had both MS and MSD recoveries (133.9% and 133.9%) above the upper limit of 125% and the results are flagged as J+.

#### Lab Report 160875

- Antimony, cobalt, lead, manganese, nickel, and selenium had low MS recoveries of 29.8%, 69.2%, 67%, 74.6%, 72.6%, and 68.2%, respectively, vs the lower limit of 75%. Antimony, arsenic, beryllium, cadmium, cobalt, lead, manganese, nickel, and selenium had low MSD recoveries of 27.2%, 71.4%, 73.6%, 74.6%, 66.6%, 63.9%, 74.4%, 71.7%, and 66.2%, respectively, vs the lower limit of 75% .Low recovery was related to dilution or matrix interference. The field sample results with low recovery are qualified “J-“and non-detect results are qualified “UJ” due to a potential low bias. Barium had both MS and MSD recoveries (133.9% and 133.9%) above the upper limit of 125% and the results are flagged as J+.

#### Lab Report 1605943

- Cyanide had a low MS recovery of 70% vs. the lower limit of 80%. Cobalt, nickel, selenium, and zinc had low MSD recoveries of 72.4%, 71.1%, 69.3%, and 70.7% vs. the lower limit of 75%. Antimony and lead had low recoveries for the MS (31.9% and 74.7%, respectively, vs. 75%) and MSD (16.9% and 71.1%, respectively, vs. 75%). The field sample results with low recovery are qualified “J-“and non-detect results are qualified “UJ” due to a potential low bias.
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#### Lab Report 1605998

- Cyanide had a low MSD recovery of 89% vs. the lower limit of 90%. The field sample result with low recovery is qualified “J-“due to a potential low bias.

#### Lab Report 1609B57

- Antimony, arsenic, lead, manganese, nickel, selenium, and zinc had low MS recoveries of 48.0%, 70.9%, 72.9%, -407%, 73%, 60.5, and 74.9%, respectively, vs. a lower limit of 75%. Antimony, manganese, and selenium had low MSD recoveries of 48.1%, -195%, and 59.2%, respectively, vs. a lower limit of 75%. The field sample results with low recovery are qualified “J-“and non-detect results are qualified “UJ” due to a potential low bias. Barium had a MSD recovery of 295%, above the upper limit of 125% and the results are flagged as J+.

#### Lab Report 1609C66

- 4-Nitrophenol had a low MSD recovery of 9.01% vs the lower limit of 15% and a RPD of 58.3% vs. a limit of 41.5%. The field sample result was non-detect and the result is qualified “UJ” due to a potential low bias.

See Table A-2 for qualified data.

### DUPLICATES

#### Field Duplicates

Field duplicates were collected at a rate as stated in the approved SWMU10 Investigation Work Plan. The RPDs between the field duplicate and its associated sample were calculated and are presented in Table A-3. The field duplicates were evaluated by the following criteria:

- If an analyte was detected at a concentration greater than five times the method reporting limit, the RPD should be less than 35 percent for soil and 25 percent for ground water samples.
  - If an analyte was detected at a concentration that is less than five times the method reporting limit, then the difference between the sample and the field duplicate should not exceed the method reporting limit.
  - Duplicate RPDs are calculated by dividing the difference of the concentrations by the average of the concentrations.
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Field duplicate RPDs were within acceptance limits except for the following soil sample:

- GRO and DRO for field sample SWMU 10-9 (4-6);
- 1,2,4-Trimethylbenzene, 1,3,5-trimethylbenzene, 1-methylnaphthalene, 2-methylnaphthalene, hexachlorobutadiene, n-butylbenzene, n-propylbenzene, sec-butylbenzene, and xylene(s) for field sample SWMU10-9 (4-6'); and
- 1-methylnaphthalene and 2-methylnaphthalene for field sample SWMU10-25 (2-2.5'); and
- Iron and zinc in field sample SWMU 10-15-GW.

See Table 3A for a field duplicate summary.

### COMPLETENESS SUMMARY

The following equation was used to calculate the technical completeness:

$$\% \text{ Technical Completeness} = \left( \frac{\text{Number of usable results}}{\text{Number of reported results}} \right) \times 100$$

The technical completeness attained for SWMU10 RCRA Investigation activities was 100 percent. The completeness results are provided in Table A-4. The analytical results for the required analytes per the approved SWMU10 Work Plan were considered usable for the intended purposes and the project DQOs have been met.

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**Table A-1**  
**Sample Identification**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID           | Lab ID       | Date Collected | Sample Type |
|---------------------|--------------|----------------|-------------|
| SWMU 10 EB01        | 1504c86-001a | 4/28/2015      | EB          |
| SWMU 10-1 (2-4')    | 1504c87-001a | 4/28/2015      | N           |
| SWMU 10-1 (4-6')    | 1504c87-002a | 4/28/2015      | N           |
| SWMU 10-1 (18-20')  | 1504c87-003a | 4/28/2015      | N           |
| SWMU 10-3 (2-4')    | 1504c87-004a | 4/28/2015      | N           |
| SWMU 10-3 (6-8')    | 1504c87-005a | 4/28/2015      | N           |
| SWMU 10-3 (18-20')  | 1504c87-006a | 4/28/2015      | N           |
| SWMU 10-1-GW        | 1505001-001a | 4/28/2015      | GW          |
| Trip Blank          | 1505001-002a | NA             | TB          |
| SWMU 10 EB02        | 1505002-001a | 4/29/2015      | EB          |
| SWMU 10-5 (0-2')    | 1505003-001a | 4/29/2015      | N           |
| SWMU 10-5 (2-4')    | 1505003-002a | 4/29/2015      | N           |
| SWMU 10-5 (4-6')    | 1505003-003a | 4/29/2015      | N           |
| SWMU 10-5 (14-16')  | 1505003-004a | 4/29/2015      | N           |
| SWMU 10-5 (22-24')  | 1505003-005A | 4/29/2015      | N           |
| SWMU 10-4 (0-2')    | 1505004-001A | 4/29/2015      | N           |
| SWMU 10-4 (2-4')    | 1505004-002a | 4/29/2015      | N           |
| SWMU 10-4 (6-8')    | 1505004-003A | 4/29/2015      | N           |
| SWMU 10-4 (18-20')  | 1505004-004A | 4/29/2015      | N           |
| SWMU 10 DUP01       | 1505004-005A | 4/29/2015      | FD          |
| SWMU 10-3-GW        | 1505005-001a | 4/29/2015      | GW          |
| Trip Blank          | 1505005-002a | NA             | TB          |
| SWMU 10 EB04        | 1505047-001a | 5/1/2015       | EB          |
| SWMU 10 EB03        | 1505048-001a | 4/30/2015      | EB          |
| SWMU 10-9 (2-4')    | 1505057-001A | 4/30/2015      | N           |
| SWMU 10-9 (4-6')    | 1505057-002a | 4/30/2015      | N           |
| SWMU 10-9 (18-20')  | 1505057-003A | 4/30/2015      | N           |
| SWMU 10-9DUP02      | 1505057-004a | 4/30/2015      | FD          |
| SWMU 10-8 (2-4')    | 1505057-005a | 4/30/2015      | N           |
| SWMU 10-8 (4-6')    | 1505057-006a | 4/30/2015      | N           |
| SWMU 10-8 (18-20')  | 1505057-007A | 4/30/2015      | N           |
| SWMU 10-10 (2-4')   | 1505058-001A | 4/30/2015      | N           |
| SWMU 10-10 (4-6')   | 1505058-002a | 4/30/2015      | N           |
| SWMU 10-10 (18-20') | 1505058-003A | 4/30/2015      | N           |
| SWMU 10-7 (2-4')    | 1505059-001A | 5/1/2015       | N           |
| SWMU 10-7 (4-6')    | 1505059-002A | 5/1/2015       | N           |
| SWMU 10-7 (18-20')  | 1505059-003A | 5/1/2015       | N           |
| SWMU 10-5-GW        | 1505218-001a | 5/4/2015       | GW          |
| TRIP BLANK          | 1505218-002a | NA             | TB          |
| SWMU 10 EB05        | 1505222-001a | 5/4/2015       | EB          |
| TRIP BLANK          | 1505222-002a | NA             | TB          |
| SWMU 10-6 (2-4')    | 1505223-001A | 5/4/2015       | N           |
| SWMU 10-6 (10-12')  | 1505223-002A | 5/4/2015       | N           |
| SWMU 10-2 (0-2')    | 1505223-003A | 5/4/2015       | N           |
| SWMU 10-2 (2-4')    | 1505223-004A | 5/4/2015       | N           |
| SWMU 10 DUP03       | 1505223-005A | 5/4/2015       | FD          |
| MEOH BLANK          | 1505223-006  | 5/4/2015       | FB          |
| SWMU 10-14 (6-8')   | 1505570-001a | 5/12/2015      | N           |
| SWMU 10-14 (21-23') | 1505570-002a | 5/12/2015      | N           |
| SWMU 10-11 (4-6')   | 1505570-003a | 5/12/2015      | N           |
| SWMU 10-11 (8-10')  | 1505570-004a | 5/12/2015      | N           |
| SWMU 10-11 (18-20') | 1505570-005A | 5/12/2015      | N           |
| SWMU 10-12 (6-8')   | 1505617-001a | 5/12/2015      | N           |

**Table A-1**  
**Sample Identification**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID             | Lab ID       | Date Collected | Sample Type |
|-----------------------|--------------|----------------|-------------|
| SWMU 10-12 (20-22')   | 1505617-002a | 5/12/2015      | N           |
| SWMU 10-13 (2-4')     | 1505617-003a | 5/13/2015      | N           |
| SWMU 10-13 (6-8')     | 1505617-004a | 5/13/2015      | N           |
| SWMU 10-13 (18-20')   | 1505617-005a | 5/13/2015      | N           |
| SWMU 10-15 (2-4')     | 1505617-006a | 5/13/2015      | N           |
| SWMU 10-15 (4-6')     | 1505617-007a | 5/13/2015      | N           |
| SWMU 10-15 (18-20')   | 1505617-008a | 5/13/2015      | N           |
| SWMU 10-17 (6-8')     | 1505617-009a | 5/13/2015      | N           |
| SWMU 10 DUP04         | 1505617-010a | 5/13/2015      | FD          |
| MeOH BLANK            | 1505617-011  | 5/13/2015      | FB          |
| SWMU 10 EB06          | 1505618-001a | 5/12/2015      | EB          |
| SWMU 10-11-GW         | 1505698-001a | 5/14/2015      | GW          |
| Trip Blank            | 1505698-002a | NA             | TB          |
| SWMU 10-12-GW         | 1505700-001a | 5/14/2015      | GW          |
| Trip Blank            | 1505700-002a | NA             | TB          |
| SWMU 10-14-GW         | 1505701-001a | 5/14/2015      | GW          |
| Trip Blank            | 1505701-002a | NA             | TB          |
| SWMU 10-16 (2-4')     | 1505705-001a | 5/13/2015      | N           |
| SWMU 10-16 (4-5.5')   | 1505705-002a | 5/13/2015      | N           |
| SWMU 10-16 (8-9')     | 1505705-003a | 5/13/2015      | N           |
| SWMU 10-15-GW         | 1505708-001a | 5/14/2015      | GW          |
| TRIP BLANK            | 1505708-002a | NA             | TB          |
| SWMU 10 DUP01GW       | 1505709-001a | 5/14/2015      | FD          |
| TRIP BLANK            | 1505709-002a | NA             | TB          |
| SWMU 10-16-GW         | 1505710-001a | 5/14/2015      | GW          |
| SWMU 10-18 (2-2.5')   | 1605874-001  | 5/16/2016      | N           |
| SWMU 10-18 (8-10')    | 1605874-002  | 5/16/2016      | N           |
| SWMU 10-18 (18-20')   | 1605874-003  | 5/16/2016      | N           |
| SWMU 10-19 (2-2.5')   | 1605875-005  | 5/17/2016      | N           |
| SWMU 10-19 (12-14')   | 1605875-006  | 5/17/2016      | N           |
| SWMU 10-19 (18-20')   | 1605875-007  | 5/17/2016      | N           |
| SWMU 10-20 (2-2.5')   | 1605875-008  | 5/17/2016      | N           |
| SWMU 10-20 (8-10')    | 1605875-001  | 5/17/2016      | N           |
| SWMU 10-20 (10-12')   | 1605875-002  | 5/17/2016      | N           |
| SWMU 10-20 (16-18')   | 1605875-003  | 5/17/2016      | N           |
| SWMU 10-20 (20-22')   | 1605875-004  | 5/17/2016      | N           |
| SWMU 10-21 (2-2.5')   | 1605943-001  | 5/18/2016      | N           |
| SWMU 10-21 (12-14')   | 1605943-002  | 5/18/2016      | N           |
| SWMU 10-21 (20-22')   | 1605943-003  | 5/18/2016      | N           |
| SWMU 10-22 (2-2.5')   | 1605943-004  | 5/18/2016      | N           |
| SWMU 10-22 (8-9')     | 1605943-005  | 5/18/2016      | N           |
| DUP01                 | 1605943-006  | 5/18/2016      | FD          |
| MeOH BLANK            | 1605943-007  | NA             | MB          |
| SWMU 10-23 (2-2.5')   | 1609B57-001  | 9/19/2016      | N           |
| SWMU 10-23 (15-16')   | 1609B57-002  | 9/19/2016      | N           |
| SWMU 10-24 (2-2.5')   | 1609B57-003  | 9/19/2016      | N           |
| SWMU 10-24 (6-8')     | 1609B57-004  | 9/19/2016      | N           |
| SWMU 10-24 (8-10')    | 1609B57-005  | 9/19/2016      | N           |
| SWMU 10-24 (15-16')   | 1609B57-006  | 9/19/2016      | N           |
| SWMU 10-25 (2-2.5')   | 1609B57-007  | 9/19/2016      | N           |
| SWMU 10-25 (10-12')   | 1609B57-008  | 9/19/2016      | N           |
| SWMU 10-25 (16.5-18') | 1609B57-009  | 9/19/2016      | N           |
| SWMU 10 DUP01         | 1609B57-010  | 9/19/2016      | FD          |

**Table A-1**  
**Sample Identification**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID     | Lab ID      | Date Collected | Sample Type |
|---------------|-------------|----------------|-------------|
| MeOH BLANK    | 1609B57-011 | NA             | MB          |
| EB091916      | 1609B57-012 | 9/19/2016      | EB          |
| SWMU 10-20-GW | 1605998-001 | 5/20/2016      | GW          |
| Trip Blank    | 1605998-002 | NA             | TB          |
| SWMU 10-21-GW | 1605998-003 | 5/20/2016      | GW          |
| SWMU 10-25-GW | 1609C66-001 | 9/21/2016      | GW          |
| Trip Blank    | 1609C66-002 | NA             | TB          |

**Notes:**

N = Normal field sample  
FD = Field duplicate  
FB = Field Blank  
NA = Not Applicable

TB = Trip Blank  
EB = Equipment Blank  
GW = Groundwater  
MB = Methanol Blank

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID     | Date Collected | Analyte                     | Result   | Units | Matrix | Qualifier | Comments  |
|---------------|----------------|-----------------------------|----------|-------|--------|-----------|---|
| SWMU 10 DUP01 | 4/29/2015      | 1,3,5-Trimethylbenzene      | 0.000422 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Methylene chloride          | 0.00051  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | sec-Butylbenzene            | 0.000534 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Ethylbenzene                | 0.000638 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 2-Hexanone                  | 0.000653 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,2,4-Trimethylbenzene      | 0.000972 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Benzene                     | 0.00104  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Xylenes, Total              | 0.00171  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Toluene                     | 0.00232  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Acetone                     | 0.0166   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Barium                      | 380      | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery  |
| SWMU 10 DUP01 | 4/29/2015      | 2-Butanone                  | 0.00253  | mg/kg | Soil   | J+/- J-   | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10 DUP01 | 4/29/2015      | 1,1,1,2-Tetrachloroethane   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,1,1-Trichloroethane       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,1,2,2-Tetrachloroethane   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,1,2-Trichloroethane       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,1-Dichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,1-Dichloroethene          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,1-Dichloropropene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,2,3-Trichlorobenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,2,3-Trichloropropane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,2,4-Trichlorobenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,2-Dibromo-3-chloropropane | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,2-Dibromoethane (EDB)     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,2-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,2-Dichloroethane (EDC)    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,2-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,3-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,3-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1,4-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 1-Methylnaphthalene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 2,2-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 2-Chlorotoluene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 2-Methylnaphthalene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 4-Chlorotoluene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 4-Isopropyltoluene          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | 4-Methyl-2-pentanone        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Bromobenzene                | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Bromodichloromethane        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10 DUP01 | 4/29/2015      | Bromoform                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID       | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|-----------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| SWMU 10 DUP01   | 4/29/2015      | Bromomethane                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Carbon disulfide               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Carbon tetrachloride           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Chlorobenzene                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Chloroethane                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Chloroform                     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Chloromethane                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | cis-1,2-DCE                    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | cis-1,3-Dichloropropene        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Dibromochloromethane           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Dibromomethane                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Dichlorodifluoromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Hexachlorobutadiene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Isopropylbenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Methyl tert-butyl ether (MTBE) | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Naphthalene                    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | n-Butylbenzene                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | n-Propylbenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Styrene                        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | tert-Butylbenzene              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Tetrachloroethene (PCE)        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | trans-1,2-DCE                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | trans-1,3-Dichloropropene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Trichloroethene (TCE)          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Trichlorofluoromethane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01   | 4/29/2015      | Vinyl chloride                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10 DUP01GW | 5/14/2015      | Benzene                        | 0.4    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 2-Butanone                     | 2.9    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Acetone                        | 16     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Methyl tert-butyl ether (MTBE) | 150    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,1,1,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,1,1-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,1,2,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,1,2-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,1-Dichloroethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,1-Dichloropropene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,2,3-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,2,3-Trichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,2,4-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,2,4-Trimethylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID       | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments   |
|-----------------|----------------|-----------------------------|--------|-------|--------|-----------|--|
| SWMU 10 DUP01GW | 5/14/2015      | 1,2-Dibromo-3-chloropropane | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,2-Dibromoethane (EDB)     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,2-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,2-Dichloroethane (EDC)    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,2-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,3,5-Trimethylbenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,3-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,3-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1,4-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 1-Methylnaphthalene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 2,2-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 2-Chlorotoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 2-Hexanone                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 2-Methylnaphthalene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 4-Chlorotoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 4-Isopropyltoluene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | 4-Methyl-2-pentanone        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Bromobenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Bromodichloromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Bromoform                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Bromomethane                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Carbon disulfide            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Carbon Tetrachloride        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Chlorobenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Chloroethane                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Chloroform                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Chloromethane               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | cis-1,2-DCE                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | cis-1,3-Dichloropropene     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Dibromochloromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Dibromomethane              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Dichlorodifluoromethane     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Ethylbenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Hexachlorobutadiene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Isopropylbenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Methylene Chloride          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Naphthalene                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | n-Butylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | n-Propylbenzene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | sec-Butylbenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10 DUP01GW | 5/14/2015      | Styrene                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |



**Table A-2**  
**Qualified Data**  
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**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID       | Date Collected | Analyte                    | Result | Units | Matrix | Qualifier | Comments  |
|-----------------|----------------|----------------------------|--------|-------|--------|-----------|---|
| SWMU 10 DUP01GW | 5/14/2015      | tert-Butylbenzene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10 DUP01GW | 5/14/2015      | Tetrachloroethene (PCE)    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10 DUP01GW | 5/14/2015      | Toluene                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10 DUP01GW | 5/14/2015      | trans-1,2-DCE              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10 DUP01GW | 5/14/2015      | trans-1,3-Dichloropropene  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10 DUP01GW | 5/14/2015      | Trichloroethene (TCE)      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10 DUP01GW | 5/14/2015      | Trichlorofluoromethane     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10 DUP01GW | 5/14/2015      | Vinyl chloride             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10 DUP01GW | 5/14/2015      | Xylenes, Total             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10 DUP04   | 5/13/2015      | Fluorene                   | 3.2    | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                            |
| SWMU 10 DUP04   | 5/13/2015      | Phenanthrene               | 6      | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                            |
| SWMU 10 DUP04   | 5/13/2015      | Naphthalene                | 8.8    | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                            |
| SWMU 10 DUP04   | 5/13/2015      | 2,4-Dimethylphenol         | 18     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                            |
| SWMU 10 DUP04   | 5/13/2015      | Phenol                     | 26     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                            |
| SWMU 10 DUP04   | 5/13/2015      | 1-Methylnaphthalene        | 28     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                            |
| SWMU 10 DUP04   | 5/13/2015      | 2-Methylphenol             | 29     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                            |
| SWMU 10 DUP04   | 5/13/2015      | 2-Methylnaphthalene        | 45     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                            |
| SWMU 10 DUP04   | 5/13/2015      | 3+4-Methylphenol           | 60     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                            |
| SWMU 10 EB05    | 5/4/2015       | 1,2,4-Trichlorobenzene     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 1,2-Dichlorobenzene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 1,3-Dichlorobenzene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 1,4-Dichlorobenzene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 1-Methylnaphthalene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2,4,5-Trichlorophenol      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2,4,6-Trichlorophenol      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2,4-Dichlorophenol         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2,4-Dimethylphenol         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2,4-Dinitrophenol          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2,6-Dinitrotoluene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2-Chloronaphthalene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2-Chlorophenol             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2-Methylnaphthalene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2-Methylphenol             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2-Nitroaniline             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 2-Nitrophenol              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 3,3'-Dichlorobenzidine     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 3+4-Methylphenol           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 3-Nitroaniline             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 4,6-Dinitro-2-methylphenol | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 4-Bromophenyl phenyl ether | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05    | 5/4/2015       | 4-Chloro-3-methylphenol    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID    | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments  |
|--------------|----------------|-----------------------------|--------|-------|--------|-----------|---|
| SWMU 10 EB05 | 5/4/2015       | 4-Chloroaniline             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | 4-Chlorophenyl phenyl ether | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | 4-Nitroaniline              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | 4-Nitrophenol               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Acenaphthene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Acenaphthylene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Aniline                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Anthracene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Azobenzene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Benz(a)anthracene           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Benzo(a)pyrene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Benzo(b)fluoranthene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Benzo(g,h,i)perylene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Benzo(k)fluoranthene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Benzoic acid                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Benzyl alcohol              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Bis(2-chloroethoxy)methane  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Bis(2-chloroethyl)ether     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Bis(2-chloroisopropyl)ether | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Bis(2-ethylhexyl)phthalate  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Butyl benzyl phthalate      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Carbazole                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Chrysene                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Dibenz(a,h)anthracene       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Dibenzofuran                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Diethyl phthalate           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Dimethyl phthalate          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Di-n-butyl phthalate        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Di-n-octyl phthalate        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Fluoranthene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Fluorene                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Hexachlorobenzene           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Hexachlorobutadiene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Hexachlorocyclopentadiene   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Hexachloroethane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Indeno(1,2,3-cd)pyrene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Isophorone                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Naphthalene                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | Nitrobenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | N-Nitrosodi-n-propylamine   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10 EB05 | 5/4/2015       | N-Nitrosodiphenylamine      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID    | Date Collected | Analyte                     | Result  | Units | Matrix | Qualifier | Comments   |
|--------------|----------------|-----------------------------|---------|-------|--------|-----------|--|
| SWMU 10 EB05 | 5/4/2015       | Phenanthrene                | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                       |
| SWMU 10 EB05 | 5/4/2015       | Phenol                      | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                       |
| SWMU 10 EB05 | 5/4/2015       | Pyrene                      | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                       |
| SWMU 10 EB05 | 5/4/2015       | Pyridine                    | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                       |
| SWMU 10 EB05 | 5/4/2015       | 2,4-Dinitrotoluene          | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                       |
| SWMU 10 EB05 | 5/4/2015       | N-Nitrosodimethylamine      | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                       |
| SWMU 10 EB05 | 5/4/2015       | Pentachlorophenol           | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                       |
| SWMU 10 EB06 | 5/12/2015      | Mercury                     | 0.00015 | mg/l  | Water  | J+        | Qualified high bias since detected in method blank & field concentration is $\leq 5\times$ blank concentration |
| SWMU 10 EB06 | 5/12/2015      | 1,2,4-Trichlorobenzene      | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 1,2-Dichlorobenzene         | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 1,3-Dichlorobenzene         | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 1,4-Dichlorobenzene         | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 1-Methylnaphthalene         | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2,4,5-Trichlorophenol       | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2,4,6-Trichlorophenol       | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2,4-Dichlorophenol          | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2,4-Dimethylphenol          | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2,4-Dinitrophenol           | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2,4-Dinitrotoluene          | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2,6-Dinitrotoluene          | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2-Chloronaphthalene         | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2-Chlorophenol              | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2-Methylnaphthalene         | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2-Methylphenol              | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2-Nitroaniline              | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 2-Nitrophenol               | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 3,3'-Dichlorobenzidine      | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 3+4-Methylphenol            | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 3-Nitroaniline              | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 4,6-Dinitro-2-methylphenol  | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 4-Bromophenyl phenyl ether  | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 4-Chloro-3-methylphenol     | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 4-Chloroaniline             | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 4-Chlorophenyl phenyl ether | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 4-Nitroaniline              | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | 4-Nitrophenol               | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | Acenaphthene                | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | Acenaphthylene              | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | Aniline                     | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | Anthracene                  | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |
| SWMU 10 EB06 | 5/12/2015      | Azobenzene                  | ND      | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time   |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID         | Date Collected | Analyte                     | Result   | Units | Matrix | Qualifier | Comments   |
|-------------------|----------------|-----------------------------|----------|-------|--------|-----------|--|
| SWMU 10 EB06      | 5/12/2015      | Benz(a)anthracene           | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Benzo(a)pyrene              | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Benzo(b)fluoranthene        | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Benzo(g,h,i)perylene        | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Benzo(k)fluoranthene        | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Benzoic acid                | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Benzyl alcohol              | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Bis(2-chloroethoxy)methane  | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Bis(2-chloroethyl)ether     | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Bis(2-chloroisopropyl)ether | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Bis(2-ethylhexyl)phthalate  | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Butyl benzyl phthalate      | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Carbazole                   | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Chrysene                    | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Dibenz(a,h)anthracene       | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Dibenzofuran                | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Diethyl phthalate           | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Dimethyl phthalate          | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Di-n-butyl phthalate        | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Di-n-octyl phthalate        | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Fluoranthene                | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Fluorene                    | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Hexachlorobenzene           | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Hexachlorobutadiene         | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Hexachlorocyclopentadiene   | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Hexachloroethane            | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Indeno(1,2,3-cd)pyrene      | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Isophorone                  | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Naphthalene                 | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Nitrobenzene                | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | N-Nitrosodimethylamine      | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | N-Nitrosodi-n-propylamine   | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | N-Nitrosodiphenylamine      | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Pentachlorophenol           | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Phenanthrene                | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Phenol                      | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Pyrene                      | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10 EB06      | 5/12/2015      | Pyridine                    | ND       | ug/l  | Water  | UJ        | Qualified low bias since extracted 2 days after the 7 day holding time |
| SWMU 10-1 (18-20) | 4/28/2015      | Methylene chloride          | 0.000566 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10-1 (18-20) | 4/28/2015      | sec-Butylbenzene            | 0.000715 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,3,5-Trimethylbenzene      | 0.000993 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch       |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID         | Date Collected | Analyte                     | Result  | Units | Matrix | Qualifier | Comments  |
|-------------------|----------------|-----------------------------|---------|-------|--------|-----------|---|
| SWMU 10-1 (18-20) | 4/28/2015      | 1,2,4-Trimethylbenzene      | 0.00162 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Ethylbenzene                | 0.00192 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Acetone                     | 0.00324 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Xylenes, Total              | 0.00647 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Toluene                     | 0.00836 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Benzene                     | 0.0146  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 2-Butanone                  | 0.00156 | mg/kg | Soil   | J+ / J-   | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,1,1,2-Tetrachloroethane   | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,1,1-Trichloroethane       | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,1,2,2-Tetrachloroethane   | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,1,2-Trichloroethane       | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,1-Dichloroethane          | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,1-Dichloroethene          | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,1-Dichloropropene         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,2,3-Trichlorobenzene      | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,2,3-Trichloropropane      | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,2,4-Trichlorobenzene      | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,2-Dibromo-3-chloropropane | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,2-Dibromoethane (EDB)     | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,2-Dichlorobenzene         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,2-Dichloroethane (EDC)    | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,2-Dichloropropane         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,3-Dichlorobenzene         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,3-Dichloropropane         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,4-Dichlorobenzene         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1-Methylnaphthalene         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 2,2-Dichloropropane         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 2-Chlorotoluene             | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 2-Hexanone                  | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 2-Methylnaphthalene         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 4-Chlorotoluene             | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 4-Isopropyltoluene          | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 4-Methyl-2-pentanone        | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Bromobenzene                | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Bromodichloromethane        | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Bromoform                   | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Bromomethane                | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Carbon disulfide            | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Carbon tetrachloride        | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Chlorobenzene               | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
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| Sample ID         | Date Collected | Analyte                        | Result   | Units | Matrix | Qualifier | Comments  |
|-------------------|----------------|--------------------------------|----------|-------|--------|-----------|---|
| SWMU 10-1 (18-20) | 4/28/2015      | Chloroethane                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Chloroform                     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Chloromethane                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | cis-1,2-DCE                    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | cis-1,3-Dichloropropene        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Dibromochloromethane           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Dibromomethane                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Dichlorodifluoromethane        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Hexachlorobutadiene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Isopropylbenzene               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Methyl tert-butyl ether (MTBE) | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Naphthalene                    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | n-Butylbenzene                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | n-Propylbenzene                | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Styrene                        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | tert-Butylbenzene              | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Tetrachloroethene (PCE)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | trans-1,2-DCE                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | trans-1,3-Dichloropropene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Trichloroethene (TCE)          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Trichlorofluoromethane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | Vinyl chloride                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (18-20) | 4/28/2015      | 1,3,5-Trimethylbenzene         | 0.000993 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (2-4)   | 4/28/2015      | 1,2,4-Trimethylbenzene         | 0.0061   | mg/kg | Soil   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq 5\times$ blank concentration  |
| SWMU 10-1 (4-6)   | 4/28/2015      | 1,3,5-Trimethylbenzene         | 0.000318 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | sec-Butylbenzene               | 0.000593 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | 1,2,4-Trimethylbenzene         | 0.000619 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | Methylene chloride             | 0.000619 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | Naphthalene                    | 0.000877 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | Benzene                        | 0.00211  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | Toluene                        | 0.00253  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | Methyl tert-butyl ether (MTBE) | 0.00669  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | Acetone                        | 0.00926  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | Ethylbenzene                   | 0.0318   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | Xylenes, Total                 | 0.197    | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | 2-Butanone                     | 0.00257  | mg/kg | Soil   | J+/ J-    | Qualified high bias since detected in method blank & field concentration $< 5\times$ blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6)   | 4/28/2015      | 1,1,1,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | 1,1,1-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | 1,1,2,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6)   | 4/28/2015      | 1,1,2-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID        | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments   |
|------------------|----------------|-----------------------------|--------|-------|--------|-----------|--|
| SWMU 10-1 (4-6') | 4/28/2015      | 1,1-Dichloroethane          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,1-Dichloroethene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,1-Dichloropropene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,2,3-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,2,3-Trichloropropane      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,2,4-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,2-Dibromo-3-chloropropane | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,2-Dibromoethane (EDB)     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,2-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,2-Dichloroethane (EDC)    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,3-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,3-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1,4-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 1-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 2,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 2-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 2-Hexanone                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 2-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 4-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 4-Isopropyltoluene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | 4-Methyl-2-pentanone        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Bromobenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Bromodichloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Bromoform                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Bromomethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Carbon disulfide            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Carbon tetrachloride        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Chlorobenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Chloroethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Chloroform                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Chloromethane               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | cis-1,2-DCE                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | cis-1,3-Dichloropropene     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Dibromochloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Dibromomethane              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Dichlorodifluoromethane     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Hexachlorobutadiene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | Isopropylbenzene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | n-Butylbenzene              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-1 (4-6') | 4/28/2015      | n-Propylbenzene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |

**Table A-2**  
**Qualified Data**  
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| Sample ID           | Date Collected | Analyte                     | Result   | Units | Matrix | Qualifier | Comments  |
|---------------------|----------------|-----------------------------|----------|-------|--------|-----------|---|
| SWMU 10-1 (4-6')    | 4/28/2015      | Styrene                     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6')    | 4/28/2015      | tert-Butylbenzene           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6')    | 4/28/2015      | Tetrachloroethene (PCE)     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6')    | 4/28/2015      | trans-1,2-DCE               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6')    | 4/28/2015      | trans-1,3-Dichloropropene   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6')    | 4/28/2015      | Trichloroethene (TCE)       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6')    | 4/28/2015      | Trichlorofluoromethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-1 (4-6')    | 4/28/2015      | Vinyl chloride              | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | n-Propylbenzene             | 0.000372 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | sec-Butylbenzene            | 0.000688 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Naphthalene                 | 0.000725 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,3,5-Trimethylbenzene      | 0.00151  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,2,4-Trimethylbenzene      | 0.00238  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Ethylbenzene                | 0.00241  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Acetone                     | 0.00422  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Toluene                     | 0.0104   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Benzene                     | 0.018    | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 2-Butanone                  | 0.00159  | mg/kg | Soil   | J+/ J-    | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,1,1,2-Tetrachloroethane   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,1,1-Trichloroethane       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,1,2,2-Tetrachloroethane   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,1,2-Trichloroethane       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,1-Dichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,1-Dichloroethene          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,1-Dichloropropene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,2,3-Trichlorobenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,2,3-Trichloropropane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,2,4-Trichlorobenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,2-Dibromo-3-chloropropane | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,2-Dibromoethane (EDB)     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,2-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,2-Dichloroethane (EDC)    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,2-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,3-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,3-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1,4-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 1-Methylnaphthalene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 2,2-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 2-Chlorotoluene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 2-Hexanone                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |



**Table A-2**  
**Qualified Data**  
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| Sample ID           | Date Collected | Analyte                        | Result   | Units | Matrix | Qualifier | Comments  |
|---------------------|----------------|--------------------------------|----------|-------|--------|-----------|---|
| SWMU 10-10 (18-20') | 4/30/2015      | 2-Methylnaphthalene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 4-Chlorotoluene                | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 4-Isopropyltoluene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | 4-Methyl-2-pentanone           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Bromobenzene                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Bromodichloromethane           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Bromoform                      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Bromomethane                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Carbon disulfide               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Carbon tetrachloride           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Chlorobenzene                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Chloroethane                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Chloroform                     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Chloromethane                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | cis-1,2-DCE                    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | cis-1,3-Dichloropropene        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Dibromochloromethane           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Dibromomethane                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Dichlorodifluoromethane        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Hexachlorobutadiene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Isopropylbenzene               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Methyl tert-butyl ether (MTBE) | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Methylene chloride             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | n-Butylbenzene                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Styrene                        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | tert-Butylbenzene              | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Tetrachloroethene (PCE)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | trans-1,2-DCE                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | trans-1,3-Dichloropropene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Trichloroethene (TCE)          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Trichlorofluoromethane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (18-20') | 4/30/2015      | Vinyl chloride                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (2-4')   | 4/30/2015      | 2-Hexanone                     | 0.000582 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (2-4')   | 4/30/2015      | Acetone                        | 0.00792  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (2-4')   | 4/30/2015      | Toluene                        | 0.000255 | mg/kg | Soil   | J+/ J-    | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4')   | 4/30/2015      | 2-Butanone                     | 0.00161  | mg/kg | Soil   | J+/ J-    | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4')   | 4/30/2015      | Methylene chloride             | 0.00051  | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (2-4')   | 4/30/2015      | 1,1,1,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-10 (2-4')   | 4/30/2015      | 1,1,1-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
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| Sample ID         | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments   |
|-------------------|----------------|-----------------------------|--------|-------|--------|-----------|--|
| SWMU 10-10 (2-4') | 4/30/2015      | 1,1,2,2-Tetrachloroethane   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,1,2-Trichloroethane       | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,1-Dichloroethane          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,1-Dichloroethene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,1-Dichloropropene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,2,3-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,2,3-Trichloropropane      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,2,4-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,2,4-Trimethylbenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,2-Dibromo-3-chloropropane | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,2-Dibromoethane (EDB)     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,2-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,2-Dichloroethane (EDC)    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,3,5-Trimethylbenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,3-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,3-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1,4-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 1-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 2,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 2-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 2-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 4-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 4-Isopropyltoluene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | 4-Methyl-2-pentanone        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Benzene                     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Bromobenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Bromodichloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Bromoform                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Bromomethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Carbon disulfide            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Carbon tetrachloride        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Chlorobenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Chloroethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Chloroform                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Chloromethane               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | cis-1,2-DCE                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | cis-1,3-Dichloropropene     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Dibromochloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Dibromomethane              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-10 (2-4') | 4/30/2015      | Dichlorodifluoromethane     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |

**Table A-2**  
**Qualified Data**  
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| Sample ID          | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|--------------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| SWMU 10-10 (2-4')  | 4/30/2015      | Ethylbenzene                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Hexachlorobutadiene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Isopropylbenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Methyl tert-butyl ether (MTBE) | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Naphthalene                    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | n-Butylbenzene                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | n-Propylbenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | sec-Butylbenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Styrene                        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | tert-Butylbenzene              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Tetrachloroethene (PCE)        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | trans-1,2-DCE                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | trans-1,3-Dichloropropene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Trichloroethene (TCE)          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Trichlorofluoromethane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Vinyl chloride                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (2-4')  | 4/30/2015      | Xylenes, Total                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-10 (4-6')  | 4/30/2015      | Fluorene                       | 2.2    | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery   |
| SWMU 10-10 (4-6')  | 4/30/2015      | Phenanthrene                   | 4.1    | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery   |
| SWMU 10-10 (4-6')  | 4/30/2015      | Naphthalene                    | 5.9    | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery   |
| SWMU 10-10 (4-6')  | 4/30/2015      | 1-Methylnaphthalene            | 17     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery   |
| SWMU 10-10 (4-6')  | 4/30/2015      | 2-Methylnaphthalene            | 25     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery   |
| SWMU 10-11 (4-6')  | 5/12/2015      | Mercury                        | 0.05   | mg/kg | Soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-11 (8-10') | 5/12/2015      | Bromomethane                   | 0.026  | mg/kg | Soil   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq 5\times$ blank concentration |
| SWMU 10-11 (8-10') | 5/12/2015      | Mercury                        | 0.014  | mg/kg | Soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-11-GW      | 5/14/2015      | 1,1-Dichloroethane             | 0.93   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | sec-Butylbenzene               | 2.2    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | Benzene                        | 2.5    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | 4-Isopropyltoluene             | 2.7    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | n-Butylbenzene                 | 3.4    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | Toluene                        | 6      | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | Isopropylbenzene               | 8.5    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | Acetone                        | 8.8    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | n-Propylbenzene                | 11     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | 1,3,5-Trimethylbenzene         | 22     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | Ethylbenzene                   | 40     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | Naphthalene                    | 45     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | 1,2,4-Trimethylbenzene         | 64     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | 1-Methylnaphthalene            | 70     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | 2-Methylnaphthalene            | 98     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-11-GW      | 5/14/2015      | Xylenes, Total                 | 230    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID     | Date Collected | Analyte                        | Result  | Units | Matrix | Qualifier | Comments   |
|---------------|----------------|--------------------------------|---------|-------|--------|-----------|--|
| SWMU 10-11-GW | 5/14/2015      | Mercury                        | 0.00075 | mg/l  | Water  | J+        | Qualified high bias since detected in method blank & field concentration is ≤ 5x blank concentration |
| SWMU 10-11-GW | 5/14/2015      | Methyl tert-butyl ether (MTBE) | 14      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,1,1,2-Tetrachloroethane      | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,1,1-Trichloroethane          | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,1,2,2-Tetrachloroethane      | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,1,2-Trichloroethane          | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,1-Dichloroethene             | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,1-Dichloropropene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,2,3-Trichlorobenzene         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,2,3-Trichloropropane         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,2,4-Trichlorobenzene         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,2-Dibromo-3-chloropropane    | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,2-Dibromoethane (EDB)        | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,2-Dichlorobenzene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,2-Dichloroethane (EDC)       | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,2-Dichloropropane            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,3-Dichlorobenzene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,3-Dichloropropane            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 1,4-Dichlorobenzene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 2,2-Dichloropropane            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 2-Butanone                     | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 2-Chlorotoluene                | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 2-Hexanone                     | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 4-Chlorotoluene                | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | 4-Methyl-2-pentanone           | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Bromobenzene                   | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Bromodichloromethane           | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Bromoform                      | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Bromomethane                   | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Carbon disulfide               | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Carbon Tetrachloride           | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Chlorobenzene                  | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Chloroethane                   | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Chloroform                     | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Chloromethane                  | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | cis-1,2-DCE                    | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | cis-1,3-Dichloropropene        | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Dibromochloromethane           | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Dibromomethane                 | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Dichlorodifluoromethane        | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |
| SWMU 10-11-GW | 5/14/2015      | Hexachlorobutadiene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                               |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID     | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments  |
|---------------|----------------|--------------------------------|--------|-------|--------|-----------|---|
| SWMU 10-11-GW | 5/14/2015      | Methylene Chloride             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-11-GW | 5/14/2015      | Styrene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-11-GW | 5/14/2015      | tert-Butylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-11-GW | 5/14/2015      | Tetrachloroethene (PCE)        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-11-GW | 5/14/2015      | trans-1,2-DCE                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-11-GW | 5/14/2015      | trans-1,3-Dichloropropene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-11-GW | 5/14/2015      | Trichloroethene (TCE)          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-11-GW | 5/14/2015      | Trichlorofluoromethane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-11-GW | 5/14/2015      | Vinyl chloride                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | Benzene                        | 0.47   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | Naphthalene                    | 0.52   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 2-Methylnaphthalene            | 1.5    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | Acetone                        | 3.4    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | Methyl tert-butyl ether (MTBE) | 13     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,2,4-Trimethylbenzene         | 0.66   | ug/l  | Water  | J+/J+/J-  | Qualified high bias since detected in method blank & trip blank and field concentration <5x blank concentration/Qualified low bias due to Method SW8270C MSD recovery and RPD for batch |
| SWMU 10-12-GW | 5/14/2015      | 1,1,1,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,1,1-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,1,2,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,1,2-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,1-Dichloroethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,1-Dichloroethene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,1-Dichloropropene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,2,3-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,2,3-Trichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,2,4-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,2-Dibromo-3-chloropropane    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,2-Dibromomethane (EDB)       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,2-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,2-Dichloroethane (EDC)       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,2-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,3,5-Trimethylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,3-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,3-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1,4-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 1-Methylnaphthalene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 2,2-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 2-Butanone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 2-Chlorotoluene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 2-Hexanone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |
| SWMU 10-12-GW | 5/14/2015      | 4-Chlorotoluene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch  |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID           | Date Collected | Analyte                   | Result | Units | Matrix | Qualifier | Comments   |
|---------------------|----------------|---------------------------|--------|-------|--------|-----------|--|
| SWMU 10-12-GW       | 5/14/2015      | 4-Isopropyltoluene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | 4-Methyl-2-pentanone      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Bromobenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Bromodichloromethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Bromoform                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Bromomethane              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Carbon disulfide          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Carbon Tetrachloride      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Chlorobenzene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Chloroethane              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Chloroform                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Chloromethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | cis-1,2-DCE               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | cis-1,3-Dichloropropene   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Dibromochloromethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Dibromomethane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Dichlorodifluoromethane   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Ethylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Hexachlorobutadiene       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Isopropylbenzene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Methylene Chloride        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | n-Butylbenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | n-Propylbenzene           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | sec-Butylbenzene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Styrene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | tert-Butylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Tetrachloroethene (PCE)   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Toluene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | trans-1,2-DCE             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | trans-1,3-Dichloropropene | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Trichloroethene (TCE)     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Trichlorofluoromethane    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Vinyl chloride            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-12-GW       | 5/14/2015      | Xylenes, Total            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-13 (6-8')   | 5/13/2015      | Bromomethane              | 0.023  | mg/kg | Soil   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq 5\times$ blank concentration |
| SWMU 10-14 (21-23') | 5/12/2015      | Mercury                   | 0.0032 | mg/kg | Soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-14 (6-8')   | 5/12/2015      | Bromomethane              | 0.023  | mg/kg | Soil   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq 5\times$ blank concentration |
| SWMU 10-14 (6-8')   | 5/12/2015      | Mercury                   | 0.0046 | mg/kg | Soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-14 (6-8')   | 5/12/2015      | Antimony                  | ND     | mg/kg | Soil   | UJ        | Qualified low since Method 6010 MS/MSD recovery low  |
| SWMU 10-14-GW       | 5/14/2015      | Toluene                   | 0.3    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW       | 5/14/2015      | Benzene                   | 0.47   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID     | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|---------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| SWMU 10-14-GW | 5/14/2015      | Acetone                        | 2.7    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Methyl tert-butyl ether (MTBE) | 16     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,2,4-Trimethylbenzene         | 0.44   | ug/l  | Water  | J+/-      | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8270C MSD recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | 1,1,1,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,1,1-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,1,2,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,1,2-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,1-Dichloroethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,1-Dichloroethene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,1-Dichloropropene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,2,3-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,2,3-Trichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,2,4-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,2-Dibromo-3-chloropropane    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,2-Dibromoethane (EDB)        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,2-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,2-Dichloroethane (EDC)       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,2-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,3,5-Trimethylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,3-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,3-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1,4-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 1-Methylnaphthalene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 2,2-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 2-Butanone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 2-Chlorotoluene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 2-Hexanone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 2-Methylnaphthalene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 4-Chlorotoluene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 4-Isopropyltoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | 4-Methyl-2-pentanone           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Bromobenzene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Bromodichloromethane           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Bromoform                      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Bromomethane                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Carbon disulfide               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Carbon Tetrachloride           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Chlorobenzene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Chloroethane                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-14-GW | 5/14/2015      | Chloroform                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID     | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|---------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| SWMU 10-14-GW | 5/14/2015      | Chloromethane                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | cis-1,2-DCE                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | cis-1,3-Dichloropropene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Dibromochloromethane           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Dibromomethane                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Dichlorodifluoromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Ethylbenzene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Hexachlorobutadiene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Isopropylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Methylene Chloride             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Naphthalene                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | n-Butylbenzene                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | n-Propylbenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | sec-Butylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Styrene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | tert-Butylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Tetrachloroethene (PCE)        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | trans-1,2-DCE                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | trans-1,3-Dichloropropene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Trichloroethene (TCE)          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Trichlorofluoromethane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Vinyl chloride                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-14-GW | 5/14/2015      | Xylenes, Total                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Benzene                        | 0.51   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 2-Butanone                     | 2.9    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Acetone                        | 9.7    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Methyl tert-butyl ether (MTBE) | 150    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,1,1,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,1,1-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,1,2,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,1,2-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,1-Dichloroethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,1-Dichloropropene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,2,3-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,2,3-Trichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,2,4-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,2,4-Trimethylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,2-Dibromo-3-chloropropane    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,2-Dibromomethane (EDB)       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,2-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |



**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID     | Date Collected | Analyte                  | Result | Units | Matrix | Qualifier | Comments   |
|---------------|----------------|--------------------------|--------|-------|--------|-----------|--|
| SWMU 10-15-GW | 5/14/2015      | 1,2-Dichloroethane (EDC) | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,2-Dichloropropane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,3,5-Trimethylbenzene   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,3-Dichlorobenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,3-Dichloropropane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1,4-Dichlorobenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 1-Methylnaphthalene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 2,2-Dichloropropane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 2-Chlorotoluene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 2-Hexanone               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 2-Methylnaphthalene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 4-Chlorotoluene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 4-Isopropyltoluene       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | 4-Methyl-2-pentanone     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Bromobenzene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Bromodichloromethane     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Bromoform                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Bromomethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Carbon disulfide         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Carbon Tetrachloride     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Chlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Chloroethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Chloroform               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Chloromethane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | cis-1,2-DCE              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | cis-1,3-Dichloropropene  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Dibromochloromethane     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Dibromomethane           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Dichlorodifluoromethane  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Ethylbenzene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Hexachlorobutadiene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Isopropylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Methylene Chloride       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Naphthalene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | n-Butylbenzene           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | n-Propylbenzene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | sec-Butylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Styrene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | tert-Butylbenzene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Tetrachloroethene (PCE)  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-15-GW | 5/14/2015      | Toluene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |

**Table A-2**  
**Qualified Data**  
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**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID           | Date Collected | Analyte                        | Result  | Units | Matrix | Qualifier | Comments   |
|---------------------|----------------|--------------------------------|---------|-------|--------|-----------|--|
| SWMU 10-15-GW       | 5/14/2015      | trans-1,2-DCE                  | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-15-GW       | 5/14/2015      | trans-1,3-Dichloropropene      | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-15-GW       | 5/14/2015      | Trichloroethene (TCE)          | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-15-GW       | 5/14/2015      | Trichlorofluoromethane         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-15-GW       | 5/14/2015      | Vinyl chloride                 | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-15-GW       | 5/14/2015      | Xylenes, Total                 | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16 (2-4')   | 5/13/2015      | 2-Butanone                     | 0.00304 | mg/kg | Soil   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration  |
| SWMU 10-16 (4-5.5') | 5/13/2015      | 2-Butanone                     | 0.00146 | mg/kg | Soil   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration  |
| SWMU 10-16 (8-9')   | 5/13/2015      | 2-Butanone                     | 0.00173 | mg/kg | Soil   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration  |
| SWMU 10-16-GW       | 5/14/2015      | Benzene                        | 0.24    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | Xylenes, Total                 | 0.35    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | cis-1,2-DCE                    | 0.39    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,1-Dichloroethene             | 0.73    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | Trichloroethene (TCE)          | 0.77    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | Acetone                        | 1.6     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,1-Dichloroethane             | 3.1     | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | Methyl tert-butyl ether (MTBE) | 19      | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,2,4-Trimethylbenzene         | 0.18    | ug/l  | Water  | J+/J-     | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8270C MSD recovery and RPD for batch |
| SWMU 10-16-GW       | 5/14/2015      | 1,2-Dichloroethane (EDC)       | 0.47    | ug/l  | Water  | J+/J-     | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8270C MSD recovery and RPD for batch |
| SWMU 10-16-GW       | 5/14/2015      | Vinyl chloride                 | 0.47    | ug/l  | Water  | J+/J-     | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8270C MSD recovery and RPD for batch |
| SWMU 10-16-GW       | 5/14/2015      | 1,1,1,2-Tetrachloroethane      | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,1,1-Trichloroethane          | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,1,2,2-Tetrachloroethane      | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,1,2-Trichloroethane          | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,1-Dichloropropene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,2,3-Trichlorobenzene         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,2,3-Trichloropropane         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,2,4-Trichlorobenzene         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,2-Dibromo-3-chloropropane    | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,2-Dibromomethane (EDB)       | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,2-Dichlorobenzene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,2-Dichloropropane            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,3,5-Trimethylbenzene         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,3-Dichlorobenzene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,3-Dichloropropane            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1,4-Dichlorobenzene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 1-Methylnaphthalene            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| SWMU 10-16-GW       | 5/14/2015      | 2,2-Dichloropropane            | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |

**Table A-2**  
**Qualified Data**  
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**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID         | Date Collected | Analyte                   | Result | Units | Matrix | Qualifier | Comments   |
|-------------------|----------------|---------------------------|--------|-------|--------|-----------|--|
| SWMU 10-16-GW     | 5/14/2015      | 2-Butanone                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | 2-Chlorotoluene           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | 2-Hexanone                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | 2-Methylnaphthalene       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | 4-Chlorotoluene           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | 4-Isopropyltoluene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | 4-Methyl-2-pentanone      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Bromobenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Bromodichloromethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Bromoform                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Bromomethane              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Carbon disulfide          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Carbon Tetrachloride      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Chlorobenzene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Chloroethane              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Chloroform                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Chloromethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | cis-1,3-Dichloropropene   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Dibromochloromethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Dibromomethane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Dichlorodifluoromethane   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Ethylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Hexachlorobutadiene       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Isopropylbenzene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Methylene Chloride        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Naphthalene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | n-Butylbenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | n-Propylbenzene           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | sec-Butylbenzene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Styrene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | tert-Butylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Tetrachloroethene (PCE)   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Toluene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | trans-1,2-DCE             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | trans-1,3-Dichloropropene | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-16-GW     | 5/14/2015      | Trichlorofluoromethane    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| SWMU 10-17 (6-8') | 5/13/2015      | Fluorene                  | 7.1    | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                           |
| SWMU 10-17 (6-8') | 5/13/2015      | Phenanthrene              | 15     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                           |
| SWMU 10-17 (6-8') | 5/13/2015      | Naphthalene               | 24     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                           |
| SWMU 10-17 (6-8') | 5/13/2015      | Phenol                    | 52     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                           |
| SWMU 10-17 (6-8') | 5/13/2015      | 2-Methylphenol            | 56     | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery                           |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID          | Date Collected | Analyte                     | Result   | Units | Matrix  | Qualifier | Comments  |
|--------------------|----------------|-----------------------------|----------|-------|---------|-----------|---|
| SWMU 10-17 (6-8')  | 5/13/2015      | 1-Methylnaphthalene         | 76       | mg/kg | Soil    | J-        | Qualified low bias due to surrogate recovery  |
| SWMU 10-17 (6-8')  | 5/13/2015      | 3+4-Methylphenol            | 100      | mg/kg | Soil    | J-        | Qualified low bias due to surrogate recovery  |
| SWMU 10-17 (6-8')  | 5/13/2015      | 2-Methylnaphthalene         | 130      | mg/kg | Soil    | J-        | Qualified low bias due to surrogate recovery  |
| SWMU 10-1-GW       | 4/29/2015      | Beryllium                   | 0.00052  | mg/l  | Water   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration   |
| SWMU 10-2 (0-2')   | 5/4/2015       | 2-Butanone                  | 0.0038   | mg/kg | Unknown | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration   |
| SWMU 10-2 (2-4')   | 5/4/2015       | 2-Butanone                  | 0.00283  | mg/kg | Unknown | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration   |
| SWMU 10-3 (18-20') | 4/28/2015      | 2-Hexanone                  | 0.000748 | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | Methylene chloride          | 0.00077  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | sec-Butylbenzene            | 0.000781 | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,3,5-Trimethylbenzene      | 0.00123  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | Acetone                     | 0.00158  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,2,4-Trimethylbenzene      | 0.00184  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | Ethylbenzene                | 0.00231  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | Xylenes, Total              | 0.00707  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | Toluene                     | 0.01     | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | Benzene                     | 0.0141   | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 2-Butanone                  | 0.00151  | mg/kg | Soil    | J+/- J-   | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,1,1,2-Tetrachloroethane   | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,1,1-Trichloroethane       | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,1,2,2-Tetrachloroethane   | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,1,2-Trichloroethane       | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,1-Dichloroethane          | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,1-Dichloroethene          | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,1-Dichloropropene         | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,2,3-Trichlorobenzene      | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,2,3-Trichloropropane      | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,2,4-Trichlorobenzene      | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,2-Dibromo-3-chloropropane | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,2-Dibromoethane (EDB)     | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,2-Dichlorobenzene         | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,2-Dichloroethane (EDC)    | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,2-Dichloropropane         | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,3-Dichlorobenzene         | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,3-Dichloropropane         | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1,4-Dichlorobenzene         | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 1-Methylnaphthalene         | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 2,2-Dichloropropane         | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 2-Chlorotoluene             | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 2-Methylnaphthalene         | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20') | 4/28/2015      | 4-Chlorotoluene             | ND       | mg/kg | Soil    | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
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| Sample ID         | Date Collected | Analyte                        | Result   | Units | Matrix | Qualifier | Comments  |
|-------------------|----------------|--------------------------------|----------|-------|--------|-----------|---|
| SWMU 10-3 (18-20) | 4/28/2015      | 4-Isopropyltoluene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | 4-Methyl-2-pentanone           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Bromobenzene                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Bromodichloromethane           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Bromoform                      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Bromomethane                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Carbon disulfide               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Carbon tetrachloride           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Chlorobenzene                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Chloroethane                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Chloroform                     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Chloromethane                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | cis-1,2-DCE                    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | cis-1,3-Dichloropropene        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Dibromochloromethane           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Dibromomethane                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Dichlorodifluoromethane        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Hexachlorobutadiene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Isopropylbenzene               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Methyl tert-butyl ether (MTBE) | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Naphthalene                    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | n-Butylbenzene                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | n-Propylbenzene                | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Styrene                        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | tert-Butylbenzene              | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Tetrachloroethene (PCE)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | trans-1,2-DCE                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | trans-1,3-Dichloropropene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Trichloroethene (TCE)          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Trichlorofluoromethane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (18-20) | 4/28/2015      | Vinyl chloride                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4)   | 4/28/2015      | Ethylbenzene                   | 0.000313 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4)   | 4/28/2015      | Methylene chloride             | 0.000387 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4)   | 4/28/2015      | 2-Hexanone                     | 0.000656 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4)   | 4/28/2015      | Benzene                        | 0.000872 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4)   | 4/28/2015      | Xylenes, Total                 | 0.00108  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4)   | 4/28/2015      | Toluene                        | 0.00116  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4)   | 4/28/2015      | Acetone                        | 0.0201   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4)   | 4/28/2015      | 2-Butanone                     | 0.00273  | mg/kg | Soil   | J+/- J-   | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4)   | 4/28/2015      | 1,1,1,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID        | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments   |
|------------------|----------------|-----------------------------|--------|-------|--------|-----------|--|
| SWMU 10-3 (2-4') | 4/28/2015      | 1,1,1-Trichloroethane       | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,1,2,2-Tetrachloroethane   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,1,2-Trichloroethane       | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,1-Dichloroethane          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,1-Dichloroethene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,1-Dichloropropene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,2,3-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,2,3-Trichloropropane      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,2,4-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,2,4-Trimethylbenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,2-Dibromo-3-chloropropane | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,2-Dibromoethane (EDB)     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,2-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,2-Dichloroethane (EDC)    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,3,3-Trimethylbenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,3-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,3-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1,4-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 1-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 2,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 2-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 2-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 4-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 4-Isopropyltoluene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | 4-Methyl-2-pentanone        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Bromobenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Bromodichloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Bromoform                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Bromomethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Carbon disulfide            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Carbon tetrachloride        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Chlorobenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Chloroethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Chloroform                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Chloromethane               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | cis-1,2-DCE                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | cis-1,3-Dichloropropene     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Dibromochloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Dibromomethane              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (2-4') | 4/28/2015      | Dichlorodifluoromethane     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID        | Date Collected | Analyte                        | Result   | Units | Matrix | Qualifier | Comments  |
|------------------|----------------|--------------------------------|----------|-------|--------|-----------|---|
| SWMU 10-3 (2-4') | 4/28/2015      | Hexachlorobutadiene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | Isopropylbenzene               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | Methyl tert-butyl ether (MTBE) | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | Naphthalene                    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | n-Butylbenzene                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | n-Propylbenzene                | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | sec-Butylbenzene               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | Styrene                        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | tert-Butylbenzene              | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | Tetrachloroethene (PCE)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | trans-1,2-DCE                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | trans-1,3-Dichloropropene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | Trichloroethene (TCE)          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | Trichlorofluoromethane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (2-4') | 4/28/2015      | Vinyl chloride                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,3,5-Trimethylbenzene         | 0.000351 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | Methylene chloride             | 0.000547 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | sec-Butylbenzene               | 0.000555 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,2,4-Trimethylbenzene         | 0.000564 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 2-Hexanone                     | 0.000588 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | Ethylbenzene                   | 0.0008   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | Xylenes, Total                 | 0.00181  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | Toluene                        | 0.00301  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | Benzene                        | 0.00319  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | Acetone                        | 0.00534  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 2-Butanone                     | 0.00136  | mg/kg | Soil   | J+/- J-   | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,1,1,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,1,1-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,1,2,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,1,2-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,1-Dichloroethane             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,1-Dichloroethene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,1-Dichloropropene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,2,3-Trichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,2,3-Trichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,2,4-Trichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,2-Dibromo-3-chloropropane    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,2-Dibromoethane (EDB)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,2-Dichlorobenzene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,2-Dichloroethane (EDC)       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
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| Sample ID        | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|------------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| SWMU 10-3 (6-8') | 4/28/2015      | 1,2-Dichloropropane            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,3-Dichlorobenzene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,3-Dichloropropane            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 1,4-Dichlorobenzene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 1-Methylnaphthalene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 2,2-Dichloropropane            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 2-Chlorotoluene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 2-Methylnaphthalene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 4-Chlorotoluene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 4-Isopropyltoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | 4-Methyl-2-pentanone           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Bromobenzene                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Bromodichloromethane           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Bromoform                      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Bromomethane                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Carbon disulfide               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Carbon tetrachloride           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Chlorobenzene                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Chloroethane                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Chloroform                     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Chloromethane                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | cis-1,2-DCE                    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | cis-1,3-Dichloropropene        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Dibromochloromethane           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Dibromomethane                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Dichlorodifluoromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Hexachlorobutadiene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Isopropylbenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Methyl tert-butyl ether (MTBE) | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Naphthalene                    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | n-Butylbenzene                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | n-Propylbenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Styrene                        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | tert-Butylbenzene              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Tetrachloroethene (PCE)        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | trans-1,2-DCE                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | trans-1,3-Dichloropropene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Trichloroethene (TCE)          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Trichlorofluoromethane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3 (6-8') | 4/28/2015      | Vinyl chloride                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-3-GW     | 4/29/2015      | Gasoline Range Organics (GRO)  | 0.033  | mg/l  | Water  | J+        | Qualified high bias due to surrogate recovery                    |



**Table A-2**  
**Qualified Data**  
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| Sample ID        | Date Collected | Analyte                     | Result   | Units | Matrix | Qualifier | Comments   |
|------------------|----------------|-----------------------------|----------|-------|--------|-----------|--|
| SWMU 10-3-GW     | 4/29/2015      | Beryllium                   | 0.0004   | mg/l  | Water  | J+        | Qualified high bias since detected in method blank & field concentration is $\leq 5x$ blank concentration  |
| SWMU 10-3-GW     | 4/29/2015      | Zinc                        | 0.023    | mg/l  | Water  | J+        | Qualified high bias since detected in method blank & field concentration is $\leq 5x$ blank concentration  |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,3,5-Trimethylbenzene      | 0.000347 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | Methylene chloride          | 0.000606 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | sec-Butylbenzene            | 0.000645 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | Ethylbenzene                | 0.000703 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,2,4-Trimethylbenzene      | 0.00077  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 2-Hexanone                  | 0.000886 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | Benzene                     | 0.00118  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | Xylenes, Total              | 0.00182  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | Toluene                     | 0.00263  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | Acetone                     | 0.0237   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 2-Butanone                  | 0.00407  | mg/kg | Soil   | J+ / J-   | Qualified high bias since detected in method blank & field concentration $< 5x$ blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,1,1,2-Tetrachloroethane   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,1,1-Trichloroethane       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,1,2,2-Tetrachloroethane   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,1,2-Trichloroethane       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,1-Dichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,1-Dichloroethene          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,1-Dichloropropene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,2,3-Trichlorobenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,2,3-Trichloropropane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,2,4-Trichlorobenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,2-Dibromo-3-chloropropane | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,2-Dibromoethane (EDB)     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,2-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,2-Dichloroethane (EDC)    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,2-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,3-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,3-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1,4-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 1-Methylnaphthalene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 2,2-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 2-Chlorotoluene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 2-Methylnaphthalene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 4-Chlorotoluene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 4-Isopropyltoluene          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | 4-Methyl-2-pentanone        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | Bromobenzene                | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (0-2') | 4/29/2015      | Bromodichloromethane        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |

**Table A-2**  
**Qualified Data**  
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| Sample ID          | Date Collected | Analyte                        | Result   | Units | Matrix | Qualifier | Comments   |
|--------------------|----------------|--------------------------------|----------|-------|--------|-----------|--|
| SWMU 10-4 (0-2')   | 4/29/2015      | Bromoform                      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Bromomethane                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Carbon disulfide               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Carbon tetrachloride           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Chlorobenzene                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Chloroethane                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Chloroform                     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Chloromethane                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | cis-1,2-DCE                    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | cis-1,3-Dichloropropene        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Dibromochloromethane           | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Dibromomethane                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Dichlorodifluoromethane        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Hexachlorobutadiene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Isopropylbenzene               | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Methyl tert-butyl ether (MTBE) | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Naphthalene                    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | n-Butylbenzene                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | n-Propylbenzene                | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Styrene                        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | tert-Butylbenzene              | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Tetrachloroethene (PCE)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | trans-1,2-DCE                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | trans-1,3-Dichloropropene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Trichloroethene (TCE)          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Trichlorofluoromethane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (0-2')   | 4/29/2015      | Vinyl chloride                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | Isopropylbenzene               | 0.000482 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | n-Propylbenzene                | 0.00051  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | Methylene chloride             | 0.000612 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | sec-Butylbenzene               | 0.000714 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | Naphthalene                    | 0.000742 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | Acetone                        | 0.0018   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | 1,3,5-Trimethylbenzene         | 0.00184  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | 1,2,4-Trimethylbenzene         | 0.00277  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | Ethylbenzene                   | 0.00291  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | Xylenes, Total                 | 0.00975  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | Toluene                        | 0.0127   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | Benzene                        | 0.0204   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | 1,1,1,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20') | 4/29/2015      | 1,1,1-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |

**Table A-2**  
**Qualified Data**  
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| Sample ID         | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments   |
|-------------------|----------------|-----------------------------|--------|-------|--------|-----------|--|
| SWMU 10-4 (18-20) | 4/29/2015      | 1,1,2,2-Tetrachloroethane   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,1,2-Trichloroethane       | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,1-Dichloroethane          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,1-Dichloroethene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,1-Dichloropropene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,2,3-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,2,3-Trichloropropane      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,2,4-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,2-Dibromo-3-chloropropane | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,2-Dibromoethane (EDB)     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,2-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,2-Dichloroethane (EDC)    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,3-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,3-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1,4-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 1-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 2,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 2-Butanone                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 2-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 2-Hexanone                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 2-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 4-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 4-Isopropyltoluene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | 4-Methyl-2-pentanone        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Bromobenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Bromodichloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Bromoform                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Bromomethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Carbon disulfide            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Carbon tetrachloride        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Chlorobenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Chloroethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Chloroform                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Chloromethane               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | cis-1,2-DCE                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | cis-1,3-Dichloropropene     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Dibromochloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Dibromomethane              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Dichlorodifluoromethane     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-4 (18-20) | 4/29/2015      | Hexachlorobutadiene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |

**Table A-2**  
**Qualified Data**  
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| Sample ID         | Date Collected | Analyte                        | Result   | Units | Matrix | Qualifier | Comments   |
|-------------------|----------------|--------------------------------|----------|-------|--------|-----------|--|
| SWMU 10-4 (18-20) | 4/29/2015      | Methyl tert-butyl ether (MTBE) | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (18-20) | 4/29/2015      | n-Butylbenzene                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (18-20) | 4/29/2015      | Styrene                        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (18-20) | 4/29/2015      | tert-Butylbenzene              | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (18-20) | 4/29/2015      | Tetrachloroethene (PCE)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (18-20) | 4/29/2015      | trans-1,2-DCE                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (18-20) | 4/29/2015      | trans-1,3-Dichloropropene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (18-20) | 4/29/2015      | Trichloroethene (TCE)          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (18-20) | 4/29/2015      | Trichlorofluoromethane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (18-20) | 4/29/2015      | Vinyl chloride                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-4 (6-8)   | 4/29/2015      | 2-Butanone                     | 0.00153  | mg/kg | Soil   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq 5\times$ blank concentration |
| SWMU 10-5 (22-24) | 4/29/2015      | n-Propylbenzene                | 0.000469 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | Methylene chloride             | 0.000565 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | sec-Butylbenzene               | 0.000766 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | Naphthalene                    | 0.000785 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1-Methylnaphthalene            | 0.000843 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 2-Methylnaphthalene            | 0.000852 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,3,5-Trimethylbenzene         | 0.00179  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | Acetone                        | 0.00186  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | Ethylbenzene                   | 0.00258  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,2,4-Trimethylbenzene         | 0.00263  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | Xylenes, Total                 | 0.00849  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | Toluene                        | 0.0112   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | Benzene                        | 0.0179   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,1,1,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,1,1-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,1,2,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,1,2-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,1-Dichloroethane             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,1-Dichloroethene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,1-Dichloropropene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,2,3-Trichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,2,3-Trichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,2,4-Trichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,2-Dibromo-3-chloropropane    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,2-Dibromoethane (EDB)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,2-Dichlorobenzene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,2-Dichloroethane (EDC)       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,2-Dichloropropane            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,3-Dichlorobenzene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-5 (22-24) | 4/29/2015      | 1,3-Dichloropropane            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |

**Table A-2**  
**Qualified Data**  
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| Sample ID         | Date Collected | Analyte                        | Result  | Units | Matrix | Qualifier | Comments  |
|-------------------|----------------|--------------------------------|---------|-------|--------|-----------|---|
| SWMU 10-5 (22-24) | 4/29/2015      | 1,4-Dichlorobenzene            | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | 2,2-Dichloropropane            | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | 2-Chlorotoluene                | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | 2-Hexanone                     | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | 4-Chlorotoluene                | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | 4-Isopropyltoluene             | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | 4-Methyl-2-pentanone           | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Bromobenzene                   | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Bromodichloromethane           | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Bromoform                      | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Bromomethane                   | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Carbon disulfide               | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Carbon tetrachloride           | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Chlorobenzene                  | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Chloroethane                   | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Chloroform                     | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Chloromethane                  | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | cis-1,2-DCE                    | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | cis-1,3-Dichloropropene        | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Dibromochloromethane           | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Dibromomethane                 | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Dichlorodifluoromethane        | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Hexachlorobutadiene            | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Isopropylbenzene               | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Methyl tert-butyl ether (MTBE) | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | n-Butylbenzene                 | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Styrene                        | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | tert-Butylbenzene              | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Tetrachloroethene (PCE)        | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | trans-1,2-DCE                  | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | trans-1,3-Dichloropropene      | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Trichloroethene (TCE)          | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Trichlorofluoromethane         | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | Vinyl chloride                 | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5 (22-24) | 4/29/2015      | 2-Butanone                     | ND      | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-5-GW      | 5/4/2015       | Gasoline Range Organics (GRO)  | 0.21    | mg/l  | Water  | J+/J+     | Qualified high bias due to surrogate recovery and Method GRO MS/MSD   |
| SWMU 10-5-GW      | 5/4/2015       | Beryllium                      | 0.00042 | mg/l  | Water  | J+        | Qualified high bias since detected in method blank & field concentration is $\leq 5 \times$ blank concentration |
| SWMU 10-5-GW      | 5/4/2015       | Bis(2-chloroethyl)ether        | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch  |
| SWMU 10-5-GW      | 5/4/2015       | N-Nitrosodimethylamine         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch  |
| SWMU 10-5-GW      | 5/4/2015       | Pentachlorophenol              | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch  |
| SWMU 10-5-GW      | 5/4/2015       | 1,2,4-Trichlorobenzene         | ND      | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch  |

**Table A-2**  
**Qualified Data**  
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| Sample ID    | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments  |
|--------------|----------------|-----------------------------|--------|-------|--------|-----------|---|
| SWMU 10-5-GW | 5/4/2015       | 1,2-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 1,3-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 1,4-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 1-Methylnaphthalene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2,4,5-Trichlorophenol       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2,4,6-Trichlorophenol       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2,4-Dichlorophenol          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2,4-Dimethylphenol          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2,4-Dinitrophenol           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2,6-Dinitrotoluene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2-Chloronaphthalene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2-Chlorophenol              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2-Methylnaphthalene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2-Methylphenol              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2-Nitroaniline              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 2-Nitrophenol               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 3,3'-Dichlorobenzidine      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 3+4-Methylphenol            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 3-Nitroaniline              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 4,6-Dinitro-2-methylphenol  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 4-Bromophenyl phenyl ether  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 4-Chloro-3-methylphenol     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 4-Chloroaniline             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 4-Chlorophenyl phenyl ether | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 4-Nitroaniline              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | 4-Nitrophenol               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Acenaphthene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Acenaphthylene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Aniline                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Anthracene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Azobenzene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Benz(a)anthracene           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Benzo(a)pyrene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Benzo(b)fluoranthene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Benzo(g,h,i)perylene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Benzo(k)fluoranthene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Benzoic acid                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Benzyl alcohol              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Bis(2-chloroethoxy)methane  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Bis(2-chloroisopropyl)ether | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |
| SWMU 10-5-GW | 5/4/2015       | Bis(2-ethylhexyl)phthalate  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method 8270C MS/MSD recoveryand RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID         | Date Collected | Analyte                   | Result   | Units | Matrix  | Qualifier | Comments  |
|-------------------|----------------|---------------------------|----------|-------|---------|-----------|---|
| SWMU 10-5-GW      | 5/4/2015       | Butyl benzyl phthalate    | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Carbazole                 | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Chrysene                  | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Dibenz(a,h)anthracene     | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Dibenzofuran              | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Diethyl phthalate         | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Dimethyl phthalate        | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Di-n-butyl phthalate      | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Di-n-octyl phthalate      | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Fluoranthene              | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Fluorene                  | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Hexachlorobenzene         | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Hexachlorobutadiene       | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Hexachlorocyclopentadiene | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Hexachloroethane          | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Indeno(1,2,3-cd)pyrene    | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Isophorone                | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Naphthalene               | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Nitrobenzene              | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | N-Nitrosodi-n-propylamine | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | N-Nitrosodiphenylamine    | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Phenanthrene              | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Phenol                    | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Pyrene                    | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | Pyridine                  | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-5-GW      | 5/4/2015       | 2,4-Dinitrotoluene        | ND       | ug/l  | Water   | UJ        | Qualified low bias due to Method 8270C MS/MSD recovery and RPD for batch                                  |
| SWMU 10-6 (2-4)   | 5/4/2015       | 2-Butanone                | 0.00176  | mg/kg | Unknown | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration |
| SWMU 10-7 (2-4)   | 5/1/2015       | 2-Butanone                | 0.00273  | mg/kg | Soil    | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration |
| SWMU 10-7 (4-6)   | 5/1/2015       | 2-Butanone                | 0.00175  | mg/kg | Soil    | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration |
| SWMU 10-8 (18-20) | 4/30/2015      | n-Propylbenzene           | 0.000428 | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Isopropylbenzene          | 0.000447 | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | sec-Butylbenzene          | 0.000714 | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Naphthalene               | 0.000809 | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,3,5-Trimethylbenzene    | 0.00159  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,2,4-Trimethylbenzene    | 0.00236  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Ethylbenzene              | 0.00285  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Acetone                   | 0.00466  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Xylenes, Total            | 0.00872  | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Toluene                   | 0.0122   | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Benzene                   | 0.0178   | mg/kg | Soil    | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Lead                      | 0.27     | mg/kg | Soil    | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration |

**Table A-2**  
**Qualified Data**  
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| Sample ID         | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments   |
|-------------------|----------------|-----------------------------|--------|-------|--------|-----------|--|
| SWMU 10-8 (18-20) | 4/30/2015      | 1,1,1,2-Tetrachloroethane   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,1,1-Trichloroethane       | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,1,2,2-Tetrachloroethane   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,1,2-Trichloroethane       | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,1-Dichloroethane          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,1-Dichloroethene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,1-Dichloropropene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,2,3-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,2,3-Trichloropropane      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,2,4-Trichlorobenzene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,2-Dibromo-3-chloropropane | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,2-Dibromoethane (EDB)     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,2-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,2-Dichloroethane (EDC)    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,3-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,3-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1,4-Dichlorobenzene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 1-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 2,2-Dichloropropane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 2-Butanone                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 2-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 2-Hexanone                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 2-Methylnaphthalene         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 4-Chlorotoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 4-Isopropyltoluene          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | 4-Methyl-2-pentanone        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Bromobenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Bromodichloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Bromoform                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Bromomethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Carbon disulfide            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Carbon tetrachloride        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Chlorobenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Chloroethane                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Chloroform                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Chloromethane               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | cis-1,2-DCE                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | cis-1,3-Dichloropropene     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Dibromochloromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-8 (18-20) | 4/30/2015      | Dibromomethane              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |



**Table A-2**  
**Qualified Data**  
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| Sample ID         | Date Collected | Analyte                        | Result   | Units | Matrix | Qualifier | Comments  |
|-------------------|----------------|--------------------------------|----------|-------|--------|-----------|---|
| SWMU 10-8 (18-20) | 4/30/2015      | Dichlorodifluoromethane        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Hexachlorobutadiene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Methyl tert-butyl ether (MTBE) | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Methylene chloride             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | n-Butylbenzene                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Styrene                        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | tert-Butylbenzene              | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Tetrachloroethene (PCE)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | trans-1,2-DCE                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | trans-1,3-Dichloropropene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Trichloroethene (TCE)          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Trichlorofluoromethane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (18-20) | 4/30/2015      | Vinyl chloride                 | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-8 (2-4)   | 4/30/2015      | Fluorene                       | 8.1      | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery  |
| SWMU 10-8 (2-4)   | 4/30/2015      | Naphthalene                    | 18       | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery  |
| SWMU 10-8 (2-4)   | 4/30/2015      | 1-Methylnaphthalene            | 68       | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery  |
| SWMU 10-8 (2-4)   | 4/30/2015      | 2-Methylnaphthalene            | 89       | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery  |
| SWMU 10-8 (4-6)   | 4/30/2015      | 1,2,4-Trimethylbenzene         | 0.021    | mg/kg | Soil   | J+        | Qualified high bias since detected in method blank & field concentration is $\leq$ 5x blank concentration |
| SWMU 10-9 (18-20) | 4/30/2015      | sec-Butylbenzene               | 0.000849 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | Naphthalene                    | 0.000873 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,3,5-Trimethylbenzene         | 0.00151  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,2,4-Trimethylbenzene         | 0.00229  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | Ethylbenzene                   | 0.00242  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | Acetone                        | 0.00359  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | Xylenes, Total                 | 0.00807  | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | Toluene                        | 0.0115   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | Benzene                        | 0.0159   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,1,1,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,1,1-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,1,2,2-Tetrachloroethane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,1,2-Trichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,1-Dichloroethane             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,1-Dichloropropene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,2,3-Trichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,2,3-Trichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,2,4-Trichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,2-Dibromo-3-chloropropane    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,2-Dibromoethane (EDB)        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,2-Dichlorobenzene            | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,2-Dichloroethane (EDC)       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
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| Sample ID         | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|-------------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| SWMU 10-9 (18-20) | 4/30/2015      | 1,2-Dichloropropane            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,3-Dichlorobenzene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,3-Dichloropropane            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 1,4-Dichlorobenzene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 1-Methylnaphthalene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 2,2-Dichloropropane            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 2-Butanone                     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 2-Chlorotoluene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 2-Hexanone                     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 2-Methylnaphthalene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 4-Chlorotoluene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 4-Isopropyltoluene             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | 4-Methyl-2-pentanone           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Bromobenzene                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Bromodichloromethane           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Bromoform                      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Bromomethane                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Carbon disulfide               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Carbon tetrachloride           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Chlorobenzene                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Chloroethane                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Chloroform                     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Chloromethane                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | cis-1,2-DCE                    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | cis-1,3-Dichloropropene        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Dibromochloromethane           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Dibromomethane                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Dichlorodifluoromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Hexachlorobutadiene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Isopropylbenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Methyl tert-butyl ether (MTBE) | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Methylene chloride             | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | n-Butylbenzene                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | n-Propylbenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Styrene                        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | tert-Butylbenzene              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Tetrachloroethene (PCE)        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | trans-1,2-DCE                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | trans-1,3-Dichloropropene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Trichloroethene (TCE)          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |
| SWMU 10-9 (18-20) | 4/30/2015      | Trichlorofluoromethane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch |

**Table A-2**  
**Qualified Data**  
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| Sample ID          | Date Collected | Analyte                     | Result   | Units | Matrix | Qualifier | Comments  |
|--------------------|----------------|-----------------------------|----------|-------|--------|-----------|---|
| SWMU 10-9 (18-20') | 4/30/2015      | Vinyl chloride              | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | Methylene chloride          | 0.000438 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | Benzene                     | 0.000788 | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | Acetone                     | 0.0184   | mg/kg | Soil   | J-        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | Bis(2-ethylhexyl)phthalate  | 1.5      | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery  |
| SWMU 10-9 (2-4')   | 4/30/2015      | Mercury                     | 0.07     | mg/kg | Soil   | J+        | Qualified high bias due to Method MS/MSD recovery   |
| SWMU 10-9 (2-4')   | 4/30/2015      | Toluene                     | 0.000797 | mg/kg | Soil   | J+/ J-    | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-9 (2-4')   | 4/30/2015      | 2-Butanone                  | 0.00234  | mg/kg | Soil   | J+/ J-    | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8260B MS/MSD recovery for batch |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,1,1,2-Tetrachloroethane   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,1,1-Trichloroethane       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,1,2,2-Tetrachloroethane   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,1,2-Trichloroethane       | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,1-Dichloroethane          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,1-Dichloroethene          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,1-Dichloropropene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,2,3-Trichlorobenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,2,3-Trichloropropane      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,2,4-Trichlorobenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,2,4-Trimethylbenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,2-Dibromo-3-chloropropane | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,2-Dibromoethane (EDB)     | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,2-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,2-Dichloroethane (EDC)    | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,2-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,3,5-Trimethylbenzene      | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,3-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,3-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1,4-Dichlorobenzene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 1-Methylnaphthalene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 2,2-Dichloropropane         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 2-Chlorotoluene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 2-Hexanone                  | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 2-Methylnaphthalene         | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 4-Chlorotoluene             | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 4-Isopropyltoluene          | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | 4-Methyl-2-pentanone        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | Bromobenzene                | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | Bromodichloromethane        | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |
| SWMU 10-9 (2-4')   | 4/30/2015      | Bromoform                   | ND       | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID        | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|------------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| SWMU 10-9 (2-4') | 4/30/2015      | Bromomethane                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Carbon disulfide               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Carbon tetrachloride           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Chlorobenzene                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Chloroethane                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Chloroform                     | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Chloromethane                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | cis-1,2-DCE                    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | cis-1,3-Dichloropropene        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Dibromochloromethane           | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Dibromomethane                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Dichlorodifluoromethane        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Ethylbenzene                   | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Hexachlorobutadiene            | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Isopropylbenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Methyl tert-butyl ether (MTBE) | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Naphthalene                    | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | n-Butylbenzene                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | n-Propylbenzene                | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | sec-Butylbenzene               | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Styrene                        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | tert-Butylbenzene              | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Tetrachloroethene (PCE)        | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | trans-1,2-DCE                  | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | trans-1,3-Dichloropropene      | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Trichloroethene (TCE)          | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Trichlorofluoromethane         | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Vinyl chloride                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (2-4') | 4/30/2015      | Xylenes, Total                 | ND     | mg/kg | Soil   | UJ        | Qualified low bias due to Method 8260B MS/MSD recovery for batch   |
| SWMU 10-9 (4-6') | 4/30/2015      | 2-Methylnaphthalene            | 1      | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery   |
| SWMU 10-9 (4-6') | 4/30/2015      | 1-Methylnaphthalene            | 1.1    | mg/kg | Soil   | J-        | Qualified low bias due to surrogate recovery   |
| Trip Blank       | 5/14/2015      | Ethylbenzene                   | 0.11   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| Trip Blank       | 5/14/2015      | Naphthalene                    | 0.31   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| Trip Blank       | 5/14/2015      | Xylenes, Total                 | 0.41   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| Trip Blank       | 5/14/2015      | 1-Methylnaphthalene            | 0.71   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| Trip Blank       | 5/14/2015      | 2-Methylnaphthalene            | 1.1    | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| Trip Blank       | 5/14/2015      | Toluene                        | 0.14   | ug/l  | Water  | J-        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch   |
| Trip Blank       | 5/14/2015      | 1,2,4-Trimethylbenzene         | 0.29   | ug/l  | Water  | J+/J-     | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8270C MSD recovery and RPD for batch |
| Trip Blank       | 5/14/2015      | 1,2,4-Trimethylbenzene         | 0.18   | ug/l  | Water  | J+/J-     | Qualified high bias since detected in method blank & field concentration <5x blank concentration/Qualified low bias due to Method SW8270C MSD recovery and RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID  | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments   |
|------------|----------------|-----------------------------|--------|-------|--------|-----------|--|
| Trip Blank | 5/14/2015      | 1,1,1,2-Tetrachloroethane   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,1-Trichloroethane       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,2,2-Tetrachloroethane   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,2-Trichloroethane       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1-Dichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1-Dichloroethene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1-Dichloropropene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,3-Trichlorobenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,3-Trichloropropane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,4-Trichlorobenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dibromo-3-chloropropane | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dibromoethane (EDB)     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dichloroethane (EDC)    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,3,5-Trimethylbenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,3-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,3-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,4-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2,2-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Butanone                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Chlorotoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Hexanone                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 4-Chlorotoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 4-Isopropyltoluene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 4-Methyl-2-pentanone        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Acetone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Benzene                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromobenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromodichloromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromoform                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromomethane                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Carbon disulfide            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Carbon Tetrachloride        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chlorobenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chloroethane                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chloroform                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chloromethane               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | cis-1,2-DCE                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | cis-1,3-Dichloropropene     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Dibromochloromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID  | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| Trip Blank | 5/14/2015      | Dibromomethane                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Dichlorodifluoromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Hexachlorobutadiene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Isopropylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Methyl tert-butyl ether (MTBE) | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Methylene Chloride             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | n-Butylbenzene                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | n-Propylbenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | sec-Butylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Styrene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | tert-Butylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Tetrachloroethene (PCE)        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Toluene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | trans-1,2-DCE                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | trans-1,3-Dichloropropene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Trichloroethene (TCE)          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Trichlorofluoromethane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Vinyl chloride                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,1,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,1-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,2,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,2-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1-Dichloroethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1-Dichloroethene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1-Dichloropropene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,3-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,3-Trichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,4-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dibromo-3-chloropropane    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dibromomethane (EDB)       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dichloroethane (EDC)       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,3,5-Trimethylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,3-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,3-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,4-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1-Methylnaphthalene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2,2-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Butanone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Chlorotoluene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID  | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| Trip Blank | 5/14/2015      | 2-Hexanone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Methylnaphthalene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 4-Chlorotoluene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 4-Isopropyltoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 4-Methyl-2-pentanone           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Acetone                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Benzene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromobenzene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromodichloromethane           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromoform                      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromomethane                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Carbon disulfide               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Carbon Tetrachloride           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chlorobenzene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chloroethane                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chloroform                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chloromethane                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | cis-1,2-DCE                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | cis-1,3-Dichloropropene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Dibromochloromethane           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Dibromomethane                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Dichlorodifluoromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Ethylbenzene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Hexachlorobutadiene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Isopropylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Methyl tert-butyl ether (MTBE) | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Methylene Chloride             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Naphthalene                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | n-Butylbenzene                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | n-Propylbenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | sec-Butylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Styrene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | tert-Butylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Tetrachloroethene (PCE)        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | trans-1,2-DCE                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | trans-1,3-Dichloropropene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Trichloroethene (TCE)          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Trichlorofluoromethane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Vinyl chloride                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Xylenes, Total                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,1,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID  | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments   |
|------------|----------------|-----------------------------|--------|-------|--------|-----------|--|
| Trip Blank | 5/14/2015      | 1,1,1-Trichloroethane       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,2,2-Tetrachloroethane   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1,2-Trichloroethane       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1-Dichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1-Dichloroethene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,1-Dichloropropene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,3-Trichlorobenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,3-Trichloropropane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,4-Trichlorobenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2,4-Trimethylbenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dibromo-3-chloropropane | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dibromoethane (EDB)     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dichloroethane (EDC)    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,2-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,3,5-Trimethylbenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,3-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,3-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1,4-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 1-Methylnaphthalene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2,2-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Butanone                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Chlorotoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Hexanone                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 2-Methylnaphthalene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 4-Chlorotoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 4-Isopropyltoluene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | 4-Methyl-2-pentanone        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Acetone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Benzene                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromobenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromodichloromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromoform                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Bromomethane                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Carbon disulfide            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Carbon Tetrachloride        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chlorobenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chloroethane                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chloroform                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Chloromethane               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | cis-1,2-DCE                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |



**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID  | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| Trip Blank | 5/14/2015      | cis-1,3-Dichloropropene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Dibromochloromethane           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Dibromomethane                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Dichlorodifluoromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Ethylbenzene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Hexachlorobutadiene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Isopropylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Methyl tert-butyl ether (MTBE) | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Methylene Chloride             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Naphthalene                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | n-Butylbenzene                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | n-Propylbenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | sec-Butylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Styrene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | tert-Butylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Tetrachloroethene (PCE)        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Toluene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | trans-1,2-DCE                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | trans-1,3-Dichloropropene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Trichloroethene (TCE)          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Trichlorofluoromethane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Vinyl chloride                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| Trip Blank | 5/14/2015      | Xylenes, Total                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1,1,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1,1-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1,2,2-Tetrachloroethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1,2-Trichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1-Dichloroethane             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1-Dichloroethene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1-Dichloropropene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2,3-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2,3-Trichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2,4-Trichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2,4-Trimethylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dibromo-3-chloropropane    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dibromoethane (EDB)        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dichloroethane (EDC)       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,3,5-Trimethylbenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,3-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID  | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| TRIP BLANK | 5/14/2015      | 1,3-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,4-Dichlorobenzene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1-Methylnaphthalene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2,2-Dichloropropane            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2-Butanone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2-Chlorotoluene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2-Hexanone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2-Methylnaphthalene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 4-Chlorotoluene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 4-Isopropyltoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 4-Methyl-2-pentanone           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Acetone                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Benzene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Bromobenzene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Bromodichloromethane           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Bromoform                      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Bromomethane                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Carbon disulfide               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Carbon Tetrachloride           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Chlorobenzene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Chloroethane                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Chloroform                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Chloromethane                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | cis-1,2-DCE                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | cis-1,3-Dichloropropene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Dibromochloromethane           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Dibromomethane                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Dichlorodifluoromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Ethylbenzene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Hexachlorobutadiene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Isopropylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Methyl tert-butyl ether (MTBE) | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Methylene Chloride             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Naphthalene                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | n-Butylbenzene                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | n-Propylbenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | sec-Butylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Styrene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | tert-Butylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Tetrachloroethene (PCE)        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Toluene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID  | Date Collected | Analyte                     | Result | Units | Matrix | Qualifier | Comments   |
|------------|----------------|-----------------------------|--------|-------|--------|-----------|--|
| TRIP BLANK | 5/14/2015      | trans-1,2-DCE               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | trans-1,3-Dichloropropene   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Trichloroethene (TCE)       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Trichlorofluoromethane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Vinyl chloride              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Xylenes, Total              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1,1,2-Tetrachloroethane   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1,1-Trichloroethane       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1,2,2-Tetrachloroethane   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1,2-Trichloroethane       | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1-Dichloroethane          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1-Dichloroethene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,1-Dichloropropene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2,3-Trichlorobenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2,3-Trichloropropane      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2,4-Trichlorobenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2,4-Trimethylbenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dibromo-3-chloropropane | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dibromoethane (EDB)     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dichloroethane (EDC)    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,2-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,3,5-Trimethylbenzene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,3-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,3-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1,4-Dichlorobenzene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 1-Methylnaphthalene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2,2-Dichloropropane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2-Butanone                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2-Chlorotoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2-Hexanone                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 2-Methylnaphthalene         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 4-Chlorotoluene             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 4-Isopropyltoluene          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | 4-Methyl-2-pentanone        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Acetone                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Benzene                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Bromobenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Bromodichloromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Bromoform                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |
| TRIP BLANK | 5/14/2015      | Bromomethane                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID           | Date Collected | Analyte                        | Result | Units | Matrix | Qualifier | Comments   |
|---------------------|----------------|--------------------------------|--------|-------|--------|-----------|--|
| TRIP BLANK          | 5/14/2015      | Carbon disulfide               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Carbon Tetrachloride           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Chlorobenzene                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Chloroethane                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Chloroform                     | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Chloromethane                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | cis-1,2-DCE                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | cis-1,3-Dichloropropene        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Dibromochloromethane           | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Dibromomethane                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Dichlorodifluoromethane        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Ethylbenzene                   | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Hexachlorobutadiene            | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Isopropylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Methyl tert-butyl ether (MTBE) | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Methylene Chloride             | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Naphthalene                    | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | n-Butylbenzene                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | n-Propylbenzene                | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | sec-Butylbenzene               | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Styrene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | tert-Butylbenzene              | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Tetrachloroethene (PCE)        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Toluene                        | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | trans-1,2-DCE                  | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | trans-1,3-Dichloropropene      | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Trichloroethene (TCE)          | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Trichlorofluoromethane         | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Vinyl chloride                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| TRIP BLANK          | 5/14/2015      | Xylenes, Total                 | ND     | ug/l  | Water  | UJ        | Qualified low bias due to Method SW8270C MS recovery and RPD for batch                           |
| SWMU 10-18 (2-2.5') | 5/6/2016       | Di-n-butyl phthalate           | 0.14   | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-18 (8-10')  | 5/6/2016       | Di-n-butyl phthalate           | 0.089  | mg/kg | soil   | J+        | qualified as non-detect due to presence in field sample below concentration in method blank      |
| SWMU 10-18 (2-2.5') | 5/6/2016       | mercury                        | 0.0044 | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-18 (8-10')  | 5/6/2016       | mercury                        | 0.0052 | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-18 (18-20') | 5/6/2016       | mercury                        | 0.0011 | mg/kg | soil   | J+        | qualified as non-detect due to presence in field sample at conc. <5 times conc. in method blank  |
| SWMU 10-18 (2-2.5') | 5/6/2016       | manganese                      | 200    | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-18 (8-10')  | 5/6/2016       | manganese                      | 700    | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-18 (18-20') | 5/6/2016       | manganese                      | 690    | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-18 (2-2.5') | 5/6/2016       | chromium VI                    | <0.64  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (8-10')  | 5/6/2016       | chromium VI                    | <0.64  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (18-20') | 5/6/2016       | chromium VI                    | <0.64  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID           | Date Collected | Analyte                   | Result   | Units | Matrix | Qualifier | Comments   |
|---------------------|----------------|---------------------------|----------|-------|--------|-----------|--|
| SWMU 10-18 (2-2.5') | 5/6/2016       | N-Nitrosodi-n-propylamine | < 0.1165 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (8-10')  | 5/6/2016       | N-Nitrosodi-n-propylamine | < 0.1126 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (18-20') | 5/6/2016       | N-Nitrosodi-n-propylamine | < 0.1105 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (2-2.5') | 5/6/2016       | Antimony                  | 2.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (2-2.5') | 5/6/2016       | Arsenic                   | 1.8      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-18 (2-2.5') | 5/6/2016       | Barium                    | 130      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-18 (2-2.5') | 5/6/2016       | Beryllium                 | 1.1      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-18 (2-2.5') | 5/6/2016       | Cadmium                   | < 0.0757 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (2-2.5') | 5/6/2016       | cobalt                    | 4.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery   |
| SWMU 10-18 (2-2.5') | 5/6/2016       | Lead                      | 5.2      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (2-2.5') | 5/6/2016       | manganese                 | 200      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (2-2.5') | 5/6/2016       | Nickel                    | 8.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (2-2.5') | 5/6/2016       | Selenium                  | < 1.3042 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (8-10')  | 5/6/2016       | Antimony                  | < 2.8    | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (8-10')  | 5/6/2016       | Arsenic                   | 2.2      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-18 (8-10')  | 5/6/2016       | Barium                    | 170      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-18 (8-10')  | 5/6/2016       | Beryllium                 | 1.1      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-18 (8-10')  | 5/6/2016       | Cadmium                   | < 0.15   | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (8-10')  | 5/6/2016       | cobalt                    | 6.4      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery   |
| SWMU 10-18 (8-10')  | 5/6/2016       | Lead                      | 4.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (8-10')  | 5/6/2016       | manganese                 | 700      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (8-10')  | 5/6/2016       | Nickel                    | 12       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (8-10')  | 5/6/2016       | Selenium                  | < 2.6    | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (18-20') | 5/6/2016       | Antimony                  | 2.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (18-20') | 5/6/2016       | Arsenic                   | 2        | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-18 (18-20') | 5/6/2016       | Barium                    | 1000     | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-18 (18-20') | 5/6/2016       | Beryllium                 | 0.86     | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-18 (18-20') | 5/6/2016       | Cadmium                   | < 0.074  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-18 (18-20') | 5/6/2016       | cobalt                    | 8.3      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery   |
| SWMU 10-18 (18-20') | 5/6/2016       | Lead                      | 2.6      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (18-20') | 5/6/2016       | manganese                 | 690      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (18-20') | 5/6/2016       | Nickel                    | 15       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-18 (18-20') | 5/6/2016       | Selenium                  | < 1.3    | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-19 (12-14') | 5/17/2016      | 1-Methylnaphthalene       | 7.2      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-19 (18-20') | 5/17/2016      | 1-Methylnaphthalene       | 0.0014   | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-20 (8-10')  | 5/17/2016      | 1-Methylnaphthalene       | 41       | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-20 (10-12') | 5/17/2016      | 1-Methylnaphthalene       | 25       | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-20 (16-18') | 5/17/2016      | 1-Methylnaphthalene       | 8.4      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-19 (12-14') | 5/17/2016      | Methylene chloride        | 0.05     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-20 (8-10')  | 5/17/2016      | Methylene chloride        | 0.51     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-20 (10-12') | 5/17/2016      | Methylene chloride        | 0.5      | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID           | Date Collected | Analyte   | Result   | Units | Matrix | Qualifier | Comments   |
|---------------------|----------------|-----------|----------|-------|--------|-----------|--|
| SWMU 10-19 (2-2.5') | 5/17/2016      | mercury   | 0.002    | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-19 (12-14') | 5/17/2016      | mercury   | 0.0019   | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-19 (18-20') | 5/17/2016      | mercury   | 0.00081  | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-20 (2-2.5') | 5/17/2016      | mercury   | 2        | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-20 (8-10')  | 5/17/2016      | mercury   | 0.23     | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-20 (10-12') | 5/17/2016      | mercury   | 0.0041   | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-20 (16-18') | 5/17/2016      | mercury   | 0.37     | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-20 (20-22') | 5/17/2016      | mercury   | 0.0012   | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-19 (2-2.5') | 5/17/2016      | Antimony  | 4        | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (2-2.5') | 5/17/2016      | Arsenic   | 1.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-19 (2-2.5') | 5/17/2016      | Barium    | 160      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-19 (2-2.5') | 5/17/2016      | Beryllium | 1.2      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-19 (2-2.5') | 5/17/2016      | Cadmium   | < 0.0759 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-19 (2-2.5') | 5/17/2016      | cobalt    | 5.2      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery   |
| SWMU 10-19 (2-2.5') | 5/17/2016      | Lead      | 5.1      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (2-2.5') | 5/17/2016      | manganese | 280      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (2-2.5') | 5/17/2016      | Nickel    | 9.8      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (2-2.5') | 5/17/2016      | Selenium  | < 1.3064 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-19 (12-14') | 5/17/2016      | Antimony  | < 6.8559 | mg/kg | soil   | UJ        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (12-14') | 5/17/2016      | Arsenic   | < 4.2233 | mg/kg | soil   | UJ        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-19 (12-14') | 5/17/2016      | Barium    | 830      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-19 (12-14') | 5/17/2016      | Beryllium | 1.1      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-19 (12-14') | 5/17/2016      | Cadmium   | < 0.3668 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-19 (12-14') | 5/17/2016      | cobalt    | 6.5      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery   |
| SWMU 10-19 (12-14') | 5/17/2016      | Lead      | 5.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (12-14') | 5/17/2016      | manganese | 1200     | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (12-14') | 5/17/2016      | Nickel    | 15       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (12-14') | 5/17/2016      | Selenium  | < 6.3156 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-19 (18-20') | 5/17/2016      | Antimony  | 3.2      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (18-20') | 5/17/2016      | Arsenic   | 3        | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-19 (18-20') | 5/17/2016      | Barium    | 480      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |
| SWMU 10-19 (18-20') | 5/17/2016      | Beryllium | 1        | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-19 (18-20') | 5/17/2016      | Cadmium   | < 0.071  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-19 (18-20') | 5/17/2016      | cobalt    | 7.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery   |
| SWMU 10-19 (18-20') | 5/17/2016      | Lead      | 3.5      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (18-20') | 5/17/2016      | manganese | 410      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (18-20') | 5/17/2016      | Nickel    | 15       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-19 (18-20') | 5/17/2016      | Selenium  | < 1.219  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-20 (2-2.5') | 5/17/2016      | Antimony  | 2.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-20 (2-2.5') | 5/17/2016      | Arsenic   | 5.4      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-20 (2-2.5') | 5/17/2016      | Barium    | 490      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery  |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID           | Date Collected | Analyte   | Result   | Units | Matrix | Qualifier | Comments  |
|---------------------|----------------|-----------|----------|-------|--------|-----------|---|
| SWMU 10-20 (2-2.5') | 5/17/2016      | Beryllium | 0.77     | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery       |
| SWMU 10-20 (2-2.5') | 5/17/2016      | Cadmium   | < 0.076  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-20 (2-2.5') | 5/17/2016      | cobalt    | 4.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery        |
| SWMU 10-20 (2-2.5') | 5/17/2016      | Lead      | 21       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (2-2.5') | 5/17/2016      | manganese | 520      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (2-2.5') | 5/17/2016      | Nickel    | 12       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (2-2.5') | 5/17/2016      | Selenium  | < 1.3049 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-20 (8-10')  | 5/17/2016      | Antimony  | 2.8      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (8-10')  | 5/17/2016      | Arsenic   | 3.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery       |
| SWMU 10-20 (8-10')  | 5/17/2016      | Barium    | 410      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery   |
| SWMU 10-20 (8-10')  | 5/17/2016      | Beryllium | 1.2      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery       |
| SWMU 10-20 (8-10')  | 5/17/2016      | Cadmium   | < 0.079  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-20 (8-10')  | 5/17/2016      | cobalt    | 5.6      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery        |
| SWMU 10-20 (8-10')  | 5/17/2016      | Lead      | 2.8      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (8-10')  | 5/17/2016      | manganese | 370      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (8-10')  | 5/17/2016      | Nickel    | 11       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (8-10')  | 5/17/2016      | Selenium  | < 1.3674 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-20 (10-12') | 5/17/2016      | Antimony  | 3.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (10-12') | 5/17/2016      | Arsenic   | 1.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery       |
| SWMU 10-20 (10-12') | 5/17/2016      | Barium    | 250      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery   |
| SWMU 10-20 (10-12') | 5/17/2016      | Beryllium | 1.1      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery       |
| SWMU 10-20 (10-12') | 5/17/2016      | Cadmium   | < 0.081  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-20 (10-12') | 5/17/2016      | cobalt    | 5.6      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery        |
| SWMU 10-20 (10-12') | 5/17/2016      | Lead      | 6        | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (10-12') | 5/17/2016      | manganese | 190      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (10-12') | 5/17/2016      | Nickel    | 10       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (10-12') | 5/17/2016      | Selenium  | < 1.3864 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-20 (16-18') | 5/17/2016      | Antimony  | 2.5      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (16-18') | 5/17/2016      | Arsenic   | 2.2      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery       |
| SWMU 10-20 (16-18') | 5/17/2016      | Barium    | 210      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery   |
| SWMU 10-20 (16-18') | 5/17/2016      | Beryllium | 0.87     | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery       |
| SWMU 10-20 (16-18') | 5/17/2016      | Cadmium   | < 0.075  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-20 (16-18') | 5/17/2016      | cobalt    | 5.5      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery        |
| SWMU 10-20 (16-18') | 5/17/2016      | Lead      | 5.3      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (16-18') | 5/17/2016      | manganese | 750      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (16-18') | 5/17/2016      | Nickel    | 10       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (16-18') | 5/17/2016      | Selenium  | < 1.2956 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-20 (20-22') | 5/17/2016      | Antimony  | 3.3      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-20 (20-22') | 5/17/2016      | Arsenic   | 2.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery       |
| SWMU 10-20 (20-22') | 5/17/2016      | Barium    | 750      | mg/kg | soil   | J+        | Qualified high bias due to Method MS/MSD recovery   |
| SWMU 10-20 (20-22') | 5/17/2016      | Beryllium | 0.88     | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery       |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID           | Date Collected | Analyte                    | Result   | Units | Matrix | Qualifier | Comments   |
|---------------------|----------------|----------------------------|----------|-------|--------|-----------|--|
| SWMU 10-20 (20-22') | 5/17/2016      | Cadmium                    | < 0.072  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-20 (20-22') | 5/17/2016      | cobalt                     | 8.5      | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery   |
| SWMU 10-20 (20-22') | 5/17/2016      | Lead                       | 3.5      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-20 (20-22') | 5/17/2016      | manganese                  | 1100     | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-20 (20-22') | 5/17/2016      | Nickel                     | 14       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-20 (20-22') | 5/17/2016      | Selenium                   | < 1.2368 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-21 (20-22') | 5/17/2016      | cyanide                    | 0.044    | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery   |
| SWMU 10-22 (2-2.5') | 5/17/2016      | cyanide                    | 0.04     | mg/kg | soil   | J-        | Qualified low bias due to Method MS recovery   |
| SWMU 10-22 (8-9')   | 5/17/2016      | cyanide                    | < 0.039  | mg/kg | soil   | UJ        | Qualified low bias due to MS recovery for batch  |
| SWMU 10-21 (2-2.5') | 5/17/2016      | Bis(2-ethylhexyl)phthalate | 0.17     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-21 (12-14') | 5/17/2016      | Bis(2-ethylhexyl)phthalate | 0.14     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-21 (20-22') | 5/17/2016      | Bis(2-ethylhexyl)phthalate | 0.12     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-22 (2-2.5') | 5/17/2016      | Bis(2-ethylhexyl)phthalate | 0.19     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-22 (8-9')   | 5/17/2016      | Bis(2-ethylhexyl)phthalate | 0.17     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-21 (12-14') | 5/17/2016      | Di-n-butyl phthalate       | 0.11     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-22 (2-2.5') | 5/17/2016      | Di-n-butyl phthalate       | 0.12     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-22 (8-9')   | 5/17/2016      | Di-n-butyl phthalate       | 0.11     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-21 (2-2.5') | 5/17/2016      | Antimony                   | < 1.3    | mg/kg | soil   | UJ        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-21 (2-2.5') | 5/17/2016      | cobalt                     | 3.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-21 (2-2.5') | 5/17/2016      | Lead                       | 2.1      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-21 (2-2.5') | 5/17/2016      | Nickel                     | 6        | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-21 (2-2.5') | 5/17/2016      | Selenium                   | < 1.2    | mg/kg | soil   | UJ        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-21 (2-2.5') | 5/17/2016      | zinc                       | 16       | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-21 (12-14') | 5/17/2016      | Antimony                   | < 7.3205 | mg/kg | soil   | UJ        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-21 (12-14') | 5/17/2016      | cobalt                     | 5.6      | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-21 (12-14') | 5/17/2016      | Lead                       | 5.1      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-21 (12-14') | 5/17/2016      | Nickel                     | 11       | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-21 (12-14') | 5/17/2016      | Selenium                   | < 1.3487 | mg/kg | soil   | UJ        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-21 (12-14') | 5/17/2016      | zinc                       | 20       | mg/kg | soil   | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-20-GW       | 5/20/2016      | cyanide                    | 1.05     | mg/l  | Water  | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-20-GW       | 5/20/2016      | mercury                    | 0.00016  | mg/l  | Water  | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-20-GW       | 5/20/2016      | methylene chloride         | 3.4      | ug/l  | Water  | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-20-GW       | 5/20/2016      | cyanide                    | 1.05     | mg/l  | Water  | J-        | Qualified low bias due to Method MSD recovery  |
| SWMU 10-23 (2-2.5') | 9/19/2016      | Antimony                   | < 2.2222 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-23 (2-2.5') | 9/19/2016      | Arsenic                    | < 1.9626 | mg/kg | soil   | UJ        | Qualified low bias due to MS recovery for batch  |
| SWMU 10-23 (2-2.5') | 9/19/2016      | Barium                     | 220      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-23 (2-2.5') | 9/19/2016      | Lead                       | 2.6      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-23 (2-2.5') | 9/19/2016      | Nickel                     | 12       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-23 (2-2.5') | 9/19/2016      | Selenium                   | < 4.0179 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-23 (2-2.5') | 9/19/2016      | zinc                       | 21       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-23 (15-16') | 9/19/2016      | Antimony                   | < 1.1461 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |



**Table A-2**  
**Qualified Data**  
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**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID           | Date Collected | Analyte  | Result   | Units | Matrix | Qualifier | Comments  |
|---------------------|----------------|----------|----------|-------|--------|-----------|---|
| SWMU 10-23 (15-16') | 9/19/2016      | Arsenic  | 1        | mg/kg | soil   | J-        | Qualified low bias due to MS recovery for batch     |
| SWMU 10-23 (15-16') | 9/19/2016      | Barium   | 730      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-23 (15-16') | 9/19/2016      | Lead     | 0.95     | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-23 (15-16') | 9/19/2016      | Nickel   | 13       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-23 (15-16') | 9/19/2016      | Selenium | < 2.0723 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-23 (15-16') | 9/19/2016      | zinc     | 25       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (2-2.5') | 9/19/2016      | Antimony | 1.1      | mg/kg | soil   | J-        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-24 (2-2.5') | 9/19/2016      | Arsenic  | < 0.9076 | mg/kg | soil   | UJ        | Qualified low bias due to MS recovery for batch     |
| SWMU 10-24 (2-2.5') | 9/19/2016      | Barium   | 190      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (2-2.5') | 9/19/2016      | Lead     | 2.3      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (2-2.5') | 9/19/2016      | Nickel   | 5.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (2-2.5') | 9/19/2016      | Selenium | < 1.858  | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-24 (2-2.5') | 9/19/2016      | zinc     | 11       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (6-8')   | 9/19/2016      | Antimony | < 1.2345 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-24 (6-8')   | 9/19/2016      | Arsenic  | < 1.0903 | mg/kg | soil   | UJ        | Qualified low bias due to MS recovery for batch     |
| SWMU 10-24 (6-8')   | 9/19/2016      | Barium   | 120      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (6-8')   | 9/19/2016      | Lead     | 3.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (6-8')   | 9/19/2016      | Nickel   | 9.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (6-8')   | 9/19/2016      | Selenium | < 2.2321 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-24 (6-8')   | 9/19/2016      | zinc     | 18       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (8-10')  | 9/19/2016      | Antimony | < 2.4091 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-24 (8-10')  | 9/19/2016      | Arsenic  | < 2.1276 | mg/kg | soil   | UJ        | Qualified low bias due to MS recovery for batch     |
| SWMU 10-24 (8-10')  | 9/19/2016      | Barium   | 130      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (8-10')  | 9/19/2016      | Lead     | 1.5      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (8-10')  | 9/19/2016      | Nickel   | 12       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (8-10')  | 9/19/2016      | Selenium | < 4.3557 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-24 (8-10')  | 9/19/2016      | zinc     | 22       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (15-16') | 9/19/2016      | Antimony | 2.1      | mg/kg | soil   | J-        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-24 (15-16') | 9/19/2016      | Arsenic  | < 1.0181 | mg/kg | soil   | UJ        | Qualified low bias due to MS recovery for batch     |
| SWMU 10-24 (15-16') | 9/19/2016      | Barium   | 510      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (15-16') | 9/19/2016      | Lead     | 1.1      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (15-16') | 9/19/2016      | Nickel   | 16       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-24 (15-16') | 9/19/2016      | Selenium | < 2.0843 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-24 (15-16') | 9/19/2016      | zinc     | 26       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-25 (2-2.5') | 9/19/2016      | Antimony | < 1.1611 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-25 (2-2.5') | 9/19/2016      | Arsenic  | < 1.0254 | mg/kg | soil   | UJ        | Qualified low bias due to MS recovery for batch     |
| SWMU 10-25 (2-2.5') | 9/19/2016      | Barium   | 96       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-25 (2-2.5') | 9/19/2016      | Lead     | 2.1      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-25 (2-2.5') | 9/19/2016      | Nickel   | 5.7      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |
| SWMU 10-25 (2-2.5') | 9/19/2016      | Selenium | < 2.0993 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch |
| SWMU 10-25 (2-2.5') | 9/19/2016      | zinc     | 11       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery    |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID             | Date Collected | Analyte                    | Result   | Units | Matrix | Qualifier | Comments   |
|-----------------------|----------------|----------------------------|----------|-------|--------|-----------|--|
| SWMU 10-25 (10-12')   | 9/19/2016      | Antimony                   | < 2.4915 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-25 (10-12')   | 9/19/2016      | Arsenic                    | < 2.2003 | mg/kg | soil   | UJ        | Qualified low bias due to MS recovery for batch  |
| SWMU 10-25 (10-12')   | 9/19/2016      | Barium                     | 190      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-25 (10-12')   | 9/19/2016      | Lead                       | 2.9      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-25 (10-12')   | 9/19/2016      | Nickel                     | 14       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-25 (10-12')   | 9/19/2016      | Selenium                   | < 4.5047 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-25 (10-12')   | 9/19/2016      | zinc                       | 26       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-25 (16.5-18') | 9/19/2016      | Antimony                   | 1.6      | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-25 (16.5-18') | 9/19/2016      | Arsenic                    | < 1.0232 | mg/kg | soil   | UJ        | Qualified low bias due to MS recovery for batch  |
| SWMU 10-25 (16.5-18') | 9/19/2016      | Barium                     | 460      | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-25 (16.5-18') | 9/19/2016      | Lead                       | < 0.2001 | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-25 (16.5-18') | 9/19/2016      | Nickel                     | 15       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-25 (16.5-18') | 9/19/2016      | Selenium                   | < 2.0948 | mg/kg | soil   | UJ        | Qualified low bias due to MS/MSD recovery for batch  |
| SWMU 10-25 (16.5-18') | 9/19/2016      | zinc                       | 27       | mg/kg | soil   | J-        | Qualified low bias due to Method MS/MSD recovery   |
| SWMU 10-24 (6-8')     | 9/19/2016      | Toluene                    | 0.25     | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-24 (6-8')     | 9/19/2016      | Bromomethane               | 0.32     | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25 (2-2.5')   | 9/19/2016      | 1,2,4-Trimethylbenzene     | 5.8      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-23 (15-16')   | 9/19/2016      | 1-Methylnaphthalene        | 0.0004   | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-24 (15-16')   | 9/19/2016      | 1-Methylnaphthalene        | 0.0008   | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-25 (2-2.5')   | 9/19/2016      | 1-Methylnaphthalene        | 13       | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-23 (2-2.5')   | 9/19/2016      | Bis(2-ethylhexyl)phthalate | 0.15     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-23 (2-2.5')   | 9/19/2016      | Di-n-butyl phthalate       | 0.22     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-23 (15-16')   | 9/19/2016      | Bis(2-ethylhexyl)phthalate | 0.17     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-23 (15-16')   | 9/19/2016      | Di-n-butyl phthalate       | 0.25     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-24 (2-2.5')   | 9/19/2016      | Bis(2-ethylhexyl)phthalate | 0.16     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-24 (2-2.5')   | 9/19/2016      | Di-n-butyl phthalate       | 0.22     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-24 (8-10')    | 9/19/2016      | Bis(2-ethylhexyl)phthalate | 0.23     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-24 (8-10')    | 9/19/2016      | Di-n-butyl phthalate       | 0.28     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-24 (15-16')   | 9/19/2016      | Bis(2-ethylhexyl)phthalate | 0.16     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-24 (15-16')   | 9/19/2016      | Di-n-butyl phthalate       | 0.15     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-25 (2-2.5')   | 9/19/2016      | Bis(2-ethylhexyl)phthalate | 0.22     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-25 (2-2.5')   | 9/19/2016      | Di-n-butyl phthalate       | 0.35     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-25 (10-12')   | 9/19/2016      | Bis(2-ethylhexyl)phthalate | 0.16     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-25 (10-12')   | 9/19/2016      | Di-n-butyl phthalate       | 0.22     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-25 (16.5-18') | 9/19/2016      | Bis(2-ethylhexyl)phthalate | 0.17     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-25 (16.5-18') | 9/19/2016      | Di-n-butyl phthalate       | 0.35     | mg/kg | soil   | J+        | qualified as high due to presence in field sample at conc. <5 times conc. in method blank        |
| SWMU 10-23 (2-2.5')   | 9/19/2016      | manganese                  | 450      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-23 (2-2.5')   | 9/19/2016      | zinc                       | 21       | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-23 (15-16')   | 9/19/2016      | manganese                  | 850      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-23 (15-16')   | 9/19/2016      | zinc                       | 25       | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-24 (2-2.5')   | 9/19/2016      | manganese                  | 370      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |

**Table A-2**  
**Qualified Data**  
**SWMU10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

| Sample ID             | Date Collected | Analyte                    | Result  | Units | Matrix | Qualifier | Comments   |
|-----------------------|----------------|----------------------------|---------|-------|--------|-----------|--|
| SWMU 10-24 (2-2.5')   | 9/19/2016      | zinc                       | 11      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-24 (6-8')     | 9/19/2016      | manganese                  | 200     | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-24 (6-8')     | 9/19/2016      | zinc                       | 18      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-24 (8-10')    | 9/19/2016      | manganese                  | 290     | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-24 (8-10')    | 9/19/2016      | zinc                       | 22      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-24 (15-16')   | 9/19/2016      | manganese                  | 1400    | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-24 (15-16')   | 9/19/2016      | zinc                       | 26      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25 (2-2.5')   | 9/19/2016      | manganese                  | 500     | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25 (2-2.5')   | 9/19/2016      | zinc                       | 11      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25 (10-12')   | 9/19/2016      | manganese                  | 300     | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25 (10-12')   | 9/19/2016      | zinc                       | 26      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25 (16.5-18') | 9/19/2016      | manganese                  | 1100    | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25 (16.5-18') | 9/19/2016      | zinc                       | 27      | mg/kg | soil   | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25-GW         | 9/21/2016      | manganese (total)          | 2.1     | mg/l  | Water  | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25-GW         | 9/21/2016      | Bis(2-ethylhexyl)phthalate | 2.8     | ug/l  | Water  | J+        | qualified high due to presence in method blank, blank concentration <5 times field concentration |
| SWMU 10-25-GW         | 9/21/2016      | 4-Nitrophenol              | < 2.553 | ug/l  | Water  | UJ        | Qualified low bias due to MS/MSD recovery for batch  |

**Notes:**

UJ = Estimated reporting concentration

J- = Low bias

J+ = High bias

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                  | Parameter                      | SWMU 10-14 (0-2') | SWMU 10 DUP01    | RPD (%) |
|------------------|--------------------------------|-------------------|------------------|---------|
|                  |                                | Sample Result     | Duplicate Result |         |
| TPH (mg/kg-dry): | Gasoline Range Organics (GRO)  | 0.849320979 U     | 0.858017826 U    | NC      |
|                  | Diesel Range Organics (DRO)    | 160               | 150              | 1.6     |
|                  | Motor Oil Range Organics (MRO) | 330               | 310              | 1.6     |
| VOCs (ug/kg-dry) | 1,3,5-Trimethylbenzene         | 0.000347 J        | 0.000422 J       | 4.9     |
|                  | Methylene chloride             | 0.000606 J        | 0.00051 J        | 4.3     |
|                  | sec-Butylbenzene               | 0.000645 J        | 0.000534 J       | 4.7     |
|                  | Ethylbenzene                   | 0.000703 J        | 0.000638 J       | 2.4     |
|                  | 1,2,4-Trimethylbenzene         | 0.00077 J         | 0.000972 J       | 5.8     |
|                  | 2-Hexanone                     | 0.000886 J        | 0.000653 J       | 7.6     |
|                  | Benzene                        | 0.00118 J         | 0.00104 J        | 3.2     |
|                  | Xylenes, Total                 | 0.00182 J         | 0.00171          | 1.6     |
|                  | Toluene                        | 0.00263           | 0.00232          | 3.1     |
|                  | Acetone                        | 0.0237            | 0.0166           | 8.8     |
|                  | 2-Butanone                     | 0.00407 J         | 0.00253 J        | 11.7    |
|                  | 1,1,1,2-Tetrachloroethane      | 0.00031 U         | 0.00026 U        | NC      |
|                  | 1,1,1-Trichloroethane          | 0.00027 U         | 0.00023 U        | NC      |
|                  | 1,1,2,2-Tetrachloroethane      | 0.00042 U         | 0.00035 U        | NC      |
|                  | 1,1,2-Trichloroethane          | 0.00193 U         | 0.00159 U        | NC      |
|                  | 1,1-Dichloroethane             | 0.00045 U         | 0.00037 U        | NC      |
|                  | 1,1-Dichloroethene             | 0.00032 U         | 0.00027 U        | NC      |
|                  | 1,1-Dichloropropene            | 0.00035 U         | 0.00029 U        | NC      |
|                  | 1,2,3-Trichlorobenzene         | 0.00070 U         | 0.00058 U        | NC      |
|                  | 1,2,3-Trichloropropane         | 0.00068 U         | 0.00056 U        | NC      |
|                  | 1,2,4-Trichlorobenzene         | 0.00072 U         | 0.00060 U        | NC      |
|                  | 1,2-Dibromo-3-chloropropane    | 0.00057 U         | 0.00047 U        | NC      |
|                  | 1,2-Dibromoethane (EDB)        | 0.00023 U         | 0.00019 U        | NC      |
|                  | 1,2-Dichlorobenzene            | 0.00043 U         | 0.00035 U        | NC      |
|                  | 1,2-Dichloroethane (EDC)       | 0.00052 U         | 0.00043 U        | NC      |
|                  | 1,2-Dichloropropane            | 0.00041 U         | 0.00034 U        | NC      |
|                  | 1,3-Dichlorobenzene            | 0.00033 U         | 0.00027 U        | NC      |
|                  | 1,3-Dichloropropane            | 0.00030 U         | 0.00025 U        | NC      |
|                  | 1,4-Dichlorobenzene            | 0.00030 U         | 0.00025 U        | NC      |
|                  | 1-Methylnaphthalene            | 0.00076 U         | 0.00063 U        | NC      |
|                  | 2,2-Dichloropropane            | 0.00033 U         | 0.00028 U        | NC      |
|                  | 2-Chlorotoluene                | 0.00038 U         | 0.00031 U        | NC      |
|                  | 2-Methylnaphthalene            | 0.00075 U         | 0.00062 U        | NC      |
|                  | 4-Chlorotoluene                | 0.00037 U         | 0.00031 U        | NC      |
|                  | 4-Isopropyltoluene             | 0.00075 U         | 0.00062 U        | NC      |
|                  | 4-Methyl-2-pentanone           | 0.00081 U         | 0.00067 U        | NC      |
|                  | Bromobenzene                   | 0.00054 U         | 0.00044 U        | NC      |
|                  | Bromodichloromethane           | 0.00026 U         | 0.00022 U        | NC      |
|                  | Bromoform                      | 0.00050 U         | 0.00041 U        | NC      |
|                  | Bromomethane                   | 0.00114 U         | 0.00094 U        | NC      |
|                  | Carbon disulfide               | 0.00087 U         | 0.00072 U        | NC      |
|                  | Carbon tetrachloride           | 0.00035 U         | 0.00029 U        | NC      |
|                  | Chlorobenzene                  | 0.00020 U         | 0.00016 U        | NC      |
|                  | Chloroethane                   | 0.00193 U         | 0.00159 U        | NC      |
|                  | Chloroform                     | 0.00040 U         | 0.00033 U        | NC      |
|                  | Chloromethane                  | 0.00023 U         | 0.00019 U        | NC      |
|                  | cis-1,2-DCE                    | 0.00051 U         | 0.00042 U        | NC      |
|                  | cis-1,3-Dichloropropene        | 0.00193 U         | 0.00159 U        | NC      |
|                  | Dibromochloromethane           | 0.00022 U         | 0.00018 U        | NC      |
|                  | Dibromomethane                 | 0.00109 U         | 0.00090 U        | NC      |
|                  | Dichlorodifluoromethane        | 0.00079 U         | 0.00065 U        | NC      |
|                  | Hexachlorobutadiene            | 0.00039 U         | 0.00032 U        | NC      |

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                    | Parameter                      | SWMU 10-14 (0-2') | SWMU 10 DUP01    | RPD (%) |
|--------------------|--------------------------------|-------------------|------------------|---------|
|                    |                                | Sample Result     | Duplicate Result |         |
|                    | Isopropylbenzene               | 0.00045 U         | 0.00037 U        | NC      |
|                    | Methyl tert-butyl ether (MTBE) | 0.00081 U         | 0.00067 U        | NC      |
|                    | Naphthalene                    | 0.00063 U         | 0.00052 U        | NC      |
|                    | n-Butylbenzene                 | 0.00079 U         | 0.00065 U        | NC      |
|                    | n-Propylbenzene                | 0.00036 U         | 0.00030 U        | NC      |
|                    | Styrene                        | 0.00032 U         | 0.00026 U        | NC      |
|                    | tert-Butylbenzene              | 0.00030 U         | 0.00024 U        | NC      |
|                    | Tetrachloroethene (PCE)        | 0.00038 U         | 0.00032 U        | NC      |
|                    | trans-1,2-DCE                  | 0.00031 U         | 0.00026 U        | NC      |
|                    | trans-1,3-Dichloropropene      | 0.00025 U         | 0.00021 U        | NC      |
|                    | Trichloroethene (TCE)          | 0.00038 U         | 0.00032 U        | NC      |
|                    | Trichlorofluoromethane         | 0.00039 U         | 0.00032 U        | NC      |
|                    | Vinyl chloride                 | 0.00022 U         | 0.00018 U        | NC      |
| SVOCs (mg/kg-dry): | 1,2,4-Trichlorobenzene         | 0.11459 U         | 2.81548 U        | NC      |
|                    | 1,2-Dichlorobenzene            | 0.10707 U         | 2.63077 U        | NC      |
|                    | 1,3-Dichlorobenzene            | 0.09942 U         | 2.44287 U        | NC      |
|                    | 1,4-Dichlorobenzene            | 0.11708 U         | 2.87684 U        | NC      |
|                    | 1-Methylnaphthalene            | 0.10376 U         | 2.54942 U        | NC      |
|                    | 2,4,5-Trichlorophenol          | 0.13045 U         | 3.20523 U        | NC      |
|                    | 2,4,6-Trichlorophenol          | 0.12941 U         | 3.17980 U        | NC      |
|                    | 2,4-Dichlorophenol             | 0.11194 U         | 2.75040 U        | NC      |
|                    | 2,4-Dimethylphenol             | 0.08799 U         | 2.16194 U        | NC      |
|                    | 2,4-Dinitrophenol              | 0.04651 U         | 1.14282 U        | NC      |
|                    | 2,4-Dinitrotoluene             | 0.10014 U         | 2.46060 U        | NC      |
|                    | 2,6-Dinitrotoluene             | 0.12442 U         | 3.05723 U        | NC      |
|                    | 2-Chloronaphthalene            | 0.11993 U         | 2.94674 U        | NC      |
|                    | 2-Chlorophenol                 | 0.10446 U         | 2.56672 U        | NC      |
|                    | 2-Methylnaphthalene            | 0.10244 U         | 2.51697 U        | NC      |
|                    | 2-Methylphenol                 | 0.11222 U         | 2.75733 U        | NC      |
|                    | 2-Nitroaniline                 | 0.12737 U         | 3.12964 U        | NC      |
|                    | 2-Nitrophenol                  | 0.10061 U         | 2.47207 U        | NC      |
|                    | 3,3'-Dichlorobenzidine         | 0.08619 U         | 2.11789 U        | NC      |
|                    | 3+4-Methylphenol               | 0.11601 U         | 2.85059 U        | NC      |
|                    | 3-Nitroaniline                 | 0.10785 U         | 2.64997 U        | NC      |
|                    | 4,6-Dinitro-2-methylphenol     | 0.06086 U         | 1.49551 U        | NC      |
|                    | 4-Bromophenyl phenyl ether     | 0.11876 U         | 2.91810 U        | NC      |
|                    | 4-Chloro-3-methylphenol        | 0.11145 U         | 2.73851 U        | NC      |
|                    | 4-Chloroaniline                | 0.10352 U         | 2.54359 U        | NC      |
|                    | 4-Chlorophenyl phenyl ether    | 0.16668 U         | 4.09557 U        | NC      |
|                    | 4-Nitroaniline                 | 0.10054 U         | 2.47040 U        | NC      |
|                    | 4-Nitrophenol                  | 0.09644 U         | 2.36972 U        | NC      |
|                    | Acenaphthene                   | 0.13640 U         | 3.35154 U        | NC      |
|                    | Acenaphthylene                 | 0.11339 U         | 2.78620 U        | NC      |
|                    | Aniline                        | 0.09743 U         | 2.39403 U        | NC      |
|                    | Anthracene                     | 0.09508 U         | 2.33627 U        | NC      |
|                    | Azobenzene                     | 0.12548 U         | 3.08325 U        | NC      |
|                    | Benz(a)anthracene              | 0.08798 U         | 2.16169 U        | NC      |
|                    | Benzo(a)pyrene                 | 0.12562 U         | 3.08667 U        | NC      |
|                    | Benzo(b)fluoranthene           | 0.10629 U         | 2.61175 U        | NC      |
|                    | Benzo(g,h,i)perylene           | 0.12738 U         | 3.12986 U        | NC      |
|                    | Benzo(k)fluoranthene           | 0.12650 U         | 3.10822 U        | NC      |
|                    | Benzoic acid                   | 0.06378 U         | 1.56717 U        | NC      |
|                    | Benzyl alcohol                 | 0.09943 U         | 2.44304 U        | NC      |
|                    | Bis(2-chloroethoxy)methane     | 0.10684 U         | 2.62506 U        | NC      |
|                    | Bis(2-chloroethyl)ether        | 0.11060 U         | 2.71765 U        | NC      |

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                     | Parameter                   | SWMU 10-14 (0-2') | SWMU 10 DUP01    | RPD (%) |
|---------------------|-----------------------------|-------------------|------------------|---------|
|                     |                             | Sample Result     | Duplicate Result |         |
|                     | Bis(2-chloroisopropyl)ether | 0.08552 U         | 2.10120 U        | NC      |
|                     | Bis(2-ethylhexyl)phthalate  | 0.12576 U         | 3.09000 U        | NC      |
|                     | Butyl benzyl phthalate      | 0.13665 U         | 3.35771 U        | NC      |
|                     | Carbazole                   | 0.10425 U         | 2.56160 U        | NC      |
|                     | Chrysene                    | 0.11240 U         | 2.76177 U        | NC      |
|                     | Dibenz(a,h)anthracene       | 0.12127 U         | 2.97976 U        | NC      |
|                     | Dibenzofuran                | 0.11971 U         | 2.94148 U        | NC      |
|                     | Diethyl phthalate           | 0.12704 U         | 3.12139 U        | NC      |
|                     | Dimethyl phthalate          | 0.10302 U         | 2.53119 U        | NC      |
|                     | Di-n-butyl phthalate        | 0.12659 U         | 3.11051 U        | NC      |
|                     | Di-n-octyl phthalate        | 0.12461 U         | 3.06182 U        | NC      |
|                     | Fluoranthene                | 0.13583 U         | 3.33755 U        | NC      |
|                     | Fluorene                    | 0.15239 U         | 3.74428 U        | NC      |
|                     | Hexachlorobenzene           | 0.10565 U         | 2.59588 U        | NC      |
|                     | Hexachlorobutadiene         | 0.11074 U         | 2.72107 U        | NC      |
|                     | Hexachlorocyclopentadiene   | 0.07764 U         | 1.90769 U        | NC      |
|                     | Hexachloroethane            | 0.09964 U         | 2.44821 U        | NC      |
|                     | Indeno(1,2,3-cd)pyrene      | 0.12721 U         | 3.12562 U        | NC      |
|                     | Isophorone                  | 0.11934 U         | 2.93240 U        | NC      |
|                     | Naphthalene                 | 0.10651 U         | 2.61700 U        | NC      |
|                     | Nitrobenzene                | 0.11525 U         | 2.83176 U        | NC      |
|                     | N-Nitrosodi-n-propylamine   | 0.11603 U         | 2.85088 U        | NC      |
|                     | N-Nitrosodiphenylamine      | 0.10022 U         | 2.46238 U        | NC      |
|                     | Pentachlorophenol           | 0.06989 U         | 1.71719 U        | NC      |
|                     | Phenanthrene                | 0.11710 U         | 2.87735 U        | NC      |
|                     | Phenol                      | 0.10032 U         | 2.46489 U        | NC      |
|                     | Pyrene                      | 0.14447 U         | 3.54968 U        | NC      |
|                     | Pyridine                    | 0.09406 U         | 2.31115 U        | NC      |
| Metals (mg/kg-dry): | Antimony                    | 1.51 U            | 1.45 U           | NC      |
|                     | Arsenic                     | 1.4 J             | 2.7 J            | 15.9    |
|                     | Barium                      | 460               | 380              | 4.8     |
|                     | Beryllium                   | 0.71              | 0.59             | 4.6     |
|                     | Cadmium                     | 0.03 U            | 0.03 U           | NC      |
|                     | Chromium                    | 14                | 16               | 3.3     |
|                     | Cobalt                      | 3.8               | 3.4              | 2.8     |
|                     | Cyanide                     | 0.28 U            | 0.28 U           | NC      |
|                     | Hexavalent Chromium         | 2.26 U            | 2.26 U           | NC      |
|                     | Iron                        | 10000             | 9100             | 2.4     |
|                     | Lead                        | 3.8               | 3.6              | 1.4     |
|                     | Manganese                   | 440               | 490              | 2.7     |
|                     | Mercury                     | 0.034 J           | 0.035 J          | 0.7     |
|                     | Nickel                      | 6.5               | 5.6              | 3.7     |
|                     | Selenium                    | 1.76 U            | 1.69 U           | NC      |
|                     | Silver                      | 0.03 U            | 0.03 U           | NC      |
|                     | Vanadium                    | 17                | 21               | 5.3     |
|                     | Zinc                        | 29                | 24               | 4.7     |

**Notes:**

RPD = Relative percent difference; [(difference)/(average)]\* 100

NC = Not calculated; RPD values were not calculated for non-detects

ug/kg-dry = micrograms per kilogram dry

mg/kg-dry = milligrams per kilogram

Table 3A  
Field Duplicate Summary  
SWMU 10 Investigation Report  
Western Refining Southwest, Inc. - Gallup Refinery

|                  | Parameter                      | SWMU 10-9 (4-6') | SWMU 10-9 DUP02  | RPD (%) |
|------------------|--------------------------------|------------------|------------------|---------|
|                  |                                | Sample Result    | Duplicate Result |         |
| TPH (mg/kg-dry): | Gasoline Range Organics (GRO)  | 12 J             | 69               | 35.2    |
|                  | Diesel Range Organics (DRO)    | 300              | 530              | 13.9    |
|                  | Motor Oil Range Organics (MRO) | 120              | 170              | 8.6     |
| VOCs (ug/kg-dry) | 1,1,1,2-Tetrachloroethane      | 0.01186 U        | 0.02684 U        | NC      |
|                  | 1,1,1-Trichloroethane          | 0.02196 U        | 0.04971 U        | NC      |
|                  | 1,1,2,2-Tetrachloroethane      | 0.01752 U        | 0.03967 U        | NC      |
|                  | 1,1,2-Trichloroethane          | 0.01738 U        | 0.03933 U        | NC      |
|                  | 1,1-Dichloroethane             | 0.05784 U        | 0.13093 U        | NC      |
|                  | 1,1-Dichloroethene             | 0.00995 U        | 0.02253 U        | NC      |
|                  | 1,1-Dichloropropene            | 0.02153 U        | 0.04874 U        | NC      |
|                  | 1,2,3-Trichlorobenzene         | 0.03086 U        | 0.06986 U        | NC      |
|                  | 1,2,3-Trichloropropane         | 0.02426 U        | 0.05491 U        | NC      |
|                  | 1,2,4-Trichlorobenzene         | 0.02668 U        | 0.06040 U        | NC      |
|                  | 1,2,4-Trimethylbenzene         | 0.42             | 2.8              | 37.0    |
|                  | 1,2-Dibromo-3-chloropropane    | 0.02484 U        | 0.05622 U        | NC      |
|                  | 1,2-Dibromoethane (EDB)        | 0.01152 U        | 0.02607 U        | NC      |
|                  | 1,2-Dichlorobenzene            | 0.00872 U        | 0.01974 U        | NC      |
|                  | 1,2-Dichloroethane (EDC)       | 0.05784 U        | 0.13093 U        | NC      |
|                  | 1,2-Dichloropropane            | 0.00808 U        | 0.01830 U        | NC      |
|                  | 1,3,5-Trimethylbenzene         | 0.12 J           | 0.82             | 37.2    |
|                  | 1,3-Dichlorobenzene            | 0.01703 U        | 0.03854 U        | NC      |
|                  | 1,3-Dichloropropane            | 0.01991 U        | 0.04508 U        | NC      |
|                  | 1,4-Dichlorobenzene            | 0.01801 U        | 0.04077 U        | NC      |
|                  | 1-Methylnaphthalene            | 1.1              | 9.3              | 39.4    |
|                  | 2,2-Dichloropropane            | 0.02105 U        | 0.04765 U        | NC      |
|                  | 2-Butanone                     | 0.09296 U        | 0.21044 U        | NC      |
|                  | 2-Chlorotoluene                | 0.01634 U        | 0.03699 U        | NC      |
|                  | 2-Hexanone                     | 0.03556 U        | 0.08050 U        | NC      |
|                  | 2-Methylnaphthalene            | 1.5              | 13               | 39.7    |
|                  | 4-Chlorotoluene                | 0.01093 U        | 0.02475 U        | NC      |
|                  | 4-Isopropyltoluene             | 0.01028 U        | 0.21 J           | NC      |
|                  | 4-Methyl-2-pentanone           | 0.02509 U        | 0.05679 U        | NC      |
|                  | Acetone                        | 0.10782 U        | 0.24407 U        | NC      |
|                  | Benzene                        | 0.02144 U        | 0.04854 U        | NC      |
|                  | Bromobenzene                   | 0.01073 U        | 0.02430 U        | NC      |
|                  | Bromodichloromethane           | 0.01183 U        | 0.02677 U        | NC      |
|                  | Bromoform                      | 0.01162 U        | 0.02632 U        | NC      |
|                  | Bromomethane                   | 0.02791 U        | 0.06318 U        | NC      |
|                  | Carbon disulfide               | 0.10405 U        | 0.23554 U        | NC      |
|                  | Carbon tetrachloride           | 0.01408 U        | 0.03188 U        | NC      |
|                  | Chlorobenzene                  | 0.01113 U        | 0.02519 U        | NC      |
|                  | Chloroethane                   | 0.11192 U        | 0.25337 U        | NC      |
|                  | Chloroform                     | 0.01507 U        | 0.03412 U        | NC      |
|                  | Chloromethane                  | 0.01674 U        | 0.03789 U        | NC      |
|                  | cis-1,2-DCE                    | 0.01829 U        | 0.04141 U        | NC      |
|                  | cis-1,3-Dichloropropene        | 0.01003 U        | 0.02271 U        | NC      |
|                  | Dibromochloromethane           | 0.01084 U        | 0.02454 U        | NC      |
|                  | Dibromomethane                 | 0.01522 U        | 0.03446 U        | NC      |
|                  | Dichlorodifluoromethane        | 0.03637 U        | 0.08232 U        | NC      |
|                  | Ethylbenzene                   | 0.04 J           | 0.34             | 39.5    |
|                  | Hexachlorobutadiene            | 0.02008 U        | 0.04545 U        | NC      |
|                  | Isopropylbenzene               | 0.016 J          | 0.16 J           | 40.9    |
|                  | Methyl tert-butyl ether (MTBE) | 0.02045 U        | 0.04629 U        | NC      |
|                  | Methylene chloride             | 0.05784 U        | 0.13093 U        | NC      |
|                  | Naphthalene                    | 0.32             | 2.5              | 38.7    |
|                  | n-Butylbenzene                 | 0.065 J          | 0.4 J            | 36.0    |

**Table 3A**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                    | Parameter                   | SWMU 10-9 (4-6') | SWMU 10-9 DUP02  | RPD (%) |
|--------------------|-----------------------------|------------------|------------------|---------|
|                    |                             | Sample Result    | Duplicate Result |         |
|                    | n-Propylbenzene             | 0.052 J          | 0.34             | 36.7    |
|                    | sec-Butylbenzene            | 0.021 J          | 0.17 J           | 39.0    |
|                    | Styrene                     | 0.02654 U        | 0.06008 U        | NC      |
|                    | tert-Butylbenzene           | 0.01656 U        | 0.03749 U        | NC      |
|                    | Tetrachloroethene (PCE)     | 0.01253 U        | 0.02836 U        | NC      |
|                    | Toluene                     | 0.01721 U        | 0.03895 U        | NC      |
|                    | trans-1,2-DCE               | 0.00940 U        | 0.02127 U        | NC      |
|                    | trans-1,3-Dichloropropene   | 0.02603 U        | 0.05893 U        | NC      |
|                    | Trichloroethene (TCE)       | 0.01408 U        | 0.03188 U        | NC      |
|                    | Trichlorofluoromethane      | 0.04364 U        | 0.09880 U        | NC      |
|                    | Vinyl chloride              | 0.05784 U        | 0.13093 U        | NC      |
|                    | Xylenes, Total              | 0.26 J           | 1.9              | 38.0    |
| SVOCs (mg/kg-dry): | 1,2,4-Trichlorobenzene      | 1.01696 U        | 0.12038 U        | NC      |
|                    | 1,2-Dichlorobenzene         | 0.95024 U        | 0.11248 U        | NC      |
|                    | 1,3-Dichlorobenzene         | 0.88237 U        | 0.10445 U        | NC      |
|                    | 1,4-Dichlorobenzene         | 1.03912 U        | 0.12300 U        | NC      |
|                    | 1-Methylnaphthalene         | 1.1 J            | 0.7              | 11.1    |
|                    | 2,4,5-Trichlorophenol       | 1.15774 U        | 0.13704 U        | NC      |
|                    | 2,4,6-Trichlorophenol       | 1.14855 U        | 0.13595 U        | NC      |
|                    | 2,4-Dichlorophenol          | 0.99345 U        | 0.11760 U        | NC      |
|                    | 2,4-Dimethylphenol          | 0.78090 U        | 0.09244 U        | NC      |
|                    | 2,4-Dinitrophenol           | 0.41279 U        | 0.04886 U        | NC      |
|                    | 2,4-Dinitrotoluene          | 0.88877 U        | 0.10520 U        | NC      |
|                    | 2,6-Dinitrotoluene          | 1.10428 U        | 0.13071 U        | NC      |
|                    | 2-Chloronaphthalene         | 1.06437 U        | 0.12599 U        | NC      |
|                    | 2-Chlorophenol              | 0.92710 U        | 0.10974 U        | NC      |
|                    | 2-Methylnaphthalene         | 1 J              | 0.89             | 2.9     |
|                    | 2-Methylphenol              | 0.99595 U        | 0.11789 U        | NC      |
|                    | 2-Nitroaniline              | 1.13043 U        | 0.13381 U        | NC      |
|                    | 2-Nitrophenol               | 0.89292 U        | 0.10569 U        | NC      |
|                    | 3,3'-Dichlorobenzidine      | 0.76499 U        | 0.09055 U        | NC      |
|                    | 3+4-Methylphenol            | 1.02964 U        | 0.12188 U        | NC      |
|                    | 3-Nitroaniline              | 0.95717 U        | 0.11330 U        | NC      |
|                    | 4,6-Dinitro-2-methylphenol  | 0.54018 U        | 0.06394 U        | NC      |
|                    | 4-Bromophenyl phenyl ether  | 1.05402 U        | 0.12477 U        | NC      |
|                    | 4-Chloro-3-methylphenol     | 0.98915 U        | 0.11709 U        | NC      |
|                    | 4-Chloroaniline             | 0.91875 U        | 0.10875 U        | NC      |
|                    | 4-Chlorophenyl phenyl ether | 1.47933 U        | 0.17511 U        | NC      |
|                    | 4-Nitroaniline              | 0.89231 U        | 0.10562 U        | NC      |
|                    | 4-Nitrophenol               | 0.85595 U        | 0.10132 U        | NC      |
|                    | Acenaphthene                | 1.21058 U        | 0.14330 U        | NC      |
|                    | Acenaphthylene              | 1.00638 U        | 0.11913 U        | NC      |
|                    | Aniline                     | 0.86473 U        | 0.10236 U        | NC      |
|                    | Anthracene                  | 0.84386 U        | 0.09989 U        | NC      |
|                    | Azobenzene                  | 1.11368 U        | 0.13183 U        | NC      |
|                    | Benz(a)anthracene           | 0.78081 U        | 0.09242 U        | NC      |
|                    | Benzo(a)pyrene              | 1.11491 U        | 0.13197 U        | NC      |
|                    | Benzo(b)fluoranthene        | 0.94337 U        | 0.11167 U        | NC      |
|                    | Benzo(g,h,i)perylene        | 1.13051 U        | 0.13382 U        | NC      |
|                    | Benzo(k)fluoranthene        | 1.12270 U        | 0.13289 U        | NC      |
|                    | Benzoic acid                | 0.56607 U        | 0.06701 U        | NC      |
|                    | Benzyl alcohol              | 0.88243 U        | 0.10445 U        | NC      |
|                    | Bis(2-chloroethoxy)methane  | 0.94818 U        | 0.11224 U        | NC      |
|                    | Bis(2-chloroethyl)ether     | 0.98162 U        | 0.11619 U        | NC      |
|                    | Bis(2-chloroisopropyl)ether | 0.75896 U        | 0.08984 U        | NC      |
|                    | Bis(2-ethylhexyl)phthalate  | 1.11611 U        | 0.13212 U        | NC      |



**Table 3A**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                     | Parameter                 | SWMU 10-9 (4-6') | SWMU 10-9 DUP02  | RPD (%) |
|---------------------|---------------------------|------------------|------------------|---------|
|                     |                           | Sample Result    | Duplicate Result |         |
|                     | Butyl benzyl phthalate    | 1.21281 U        | 0.14356 U        | NC      |
|                     | Carbazole                 | 0.92525 U        | 0.10952 U        | NC      |
|                     | Chrysene                  | 0.99756 U        | 0.11808 U        | NC      |
|                     | Dibenz(a,h)anthracene     | 1.07629 U        | 0.12740 U        | NC      |
|                     | Dibenzofuran              | 1.06247 U        | 0.12577 U        | NC      |
|                     | Diethyl phthalate         | 1.12745 U        | 0.13346 U        | NC      |
|                     | Dimethyl phthalate        | 0.91427 U        | 0.10822 U        | NC      |
|                     | Di-n-butyl phthalate      | 1.12352 U        | 0.13299 U        | NC      |
|                     | Di-n-octyl phthalate      | 1.10594 U        | 0.13091 U        | NC      |
|                     | Fluoranthene              | 1.20553 U        | 0.14270 U        | NC      |
|                     | Fluorene                  | 1.35244 U        | 0.16009 U        | NC      |
|                     | Hexachlorobenzene         | 0.93764 U        | 0.11099 U        | NC      |
|                     | Hexachlorobutadiene       | 0.98286 U        | 0.11634 U        | NC      |
|                     | Hexachlorocyclopentadiene | 0.68906 U        | 0.08156 U        | NC      |
|                     | Hexachloroethane          | 0.88430 U        | 0.10467 U        | NC      |
|                     | Indeno(1,2,3-cd)pyrene    | 1.12898 U        | 0.13364 U        | NC      |
|                     | Isophorone                | 1.05919 U        | 0.12538 U        | NC      |
|                     | Naphthalene               | 0.94527 U        | 0.33             | NC      |
|                     | Nitrobenzene              | 1.02284 U        | 0.12107 U        | NC      |
|                     | N-Nitrosodi-n-propylamine | 1.02975 U        | 0.12189 U        | NC      |
|                     | N-Nitrosodiphenylamine    | 0.88942 U        | 0.10528 U        | NC      |
|                     | Pentachlorophenol         | 0.62025 U        | 0.07342 U        | NC      |
|                     | Phenanthrene              | 1.03930 U        | 0.12302 U        | NC      |
|                     | Phenol                    | 0.89032 U        | 0.10539 U        | NC      |
|                     | Pyrene                    | 1.28215 U        | 0.15177 U        | NC      |
|                     | Pyridine                  | 0.83479 U        | 0.09881 U        | NC      |
| Metals (mg/kg-dry): | Antimony                  | 1.56 U           | 1.54 U           | NC      |
|                     | Arsenic                   | 1.1 J            | 1.3 J            | 4.2     |
|                     | Barium                    | 250              | 190              | 6.8     |
|                     | Beryllium                 | 0.94             | 0.9              | 1.1     |
|                     | Cadmium                   | 0.033002519 U    | 0.03 U           | NC      |
|                     | Chromium                  | 14               | 8.5              | 12.2    |
|                     | Cobalt                    | 5.3              | 4.1              | 6.4     |
|                     | Cyanide                   | 0.300287306 U    | 0.30 U           | NC      |
|                     | Hexavalent Chromium       | 2.40229845 U     | 2.37 U           | NC      |
|                     | Iron                      | 16000            | 13000            | 5.2     |
|                     | Lead                      | 3.2              | 3.3              | 0.8     |
|                     | Manganese                 | 290              | 170              | 13.0    |
|                     | Mercury                   | 0.0086 J         | 0.0043 J         | 16.7    |
|                     | Nickel                    | 10               | 7.4              | 7.5     |
|                     | Selenium                  | 1.820599379 U    | 1.80 U           | NC      |
|                     | Silver                    | 0.035376801 U    | 0.03 U           | NC      |
|                     | Vanadium                  | 19               | 19               | 0.0     |
|                     | Zinc                      | 32               | 49               | 10.5    |

**Notes:**

RPD = Relative percent difference; [(difference)/(average)]\* 100

NC = Not calculated; RPD values were not calculated for non-detects

ug/kg-dry = micrograms per kilogram dry

mg/kg-dry = milligrams per kilogram

bold value = Field Duplicate RPD Outlier

Table A-3  
Field Duplicate Summary  
SWMU 10 Investigation Report  
Western Refining Southwest, Inc. - Gallup Refinery

|                  | Parameter                      | SWMU 10-6 (10-12') | SWMU 10 DUP03   | RPD (%) |
|------------------|--------------------------------|--------------------|-----------------|---------|
|                  |                                | Sample Result      | Field Duplicate |         |
| TPH (mg/kg-dry): | Gasoline Range Organics (GRO)  | 1.22 U             | 1.16 U          | NC      |
|                  | Diesel Range Organics (DRO)    | 5.32 U             | 5.44 U          | NC      |
|                  | Motor Oil Range Organics (MRO) | 48.08 U            | 49.16 U         | NC      |
| VOCs (ug/kg-dry) | 1,1,1,2-Tetrachloroethane      | 0.000407 U         | 0.000378 U      | NC      |
|                  | 1,1,1-Trichloroethane          | 0.000356 U         | 0.000331 U      | NC      |
|                  | 1,1,2,2-Tetrachloroethane      | 0.000550 U         | 0.000511 U      | NC      |
|                  | 1,1,2-Trichloroethane          | 0.002501 U         | 0.002325 U      | NC      |
|                  | 1,1-Dichloroethane             | 0.000584 U         | 0.000543 U      | NC      |
|                  | 1,1-Dichloroethene             | 0.000417 U         | 0.000388 U      | NC      |
|                  | 1,1-Dichloropropene            | 0.000458 U         | 0.000426 U      | NC      |
|                  | 1,2,3-Trichlorobenzene         | 0.000915 U         | 0.000851 U      | NC      |
|                  | 1,2,3-Trichloropropane         | 0.000878 U         | 0.000816 U      | NC      |
|                  | 1,2,4-Trichlorobenzene         | 0.000937 U         | 0.000871 U      | NC      |
|                  | 1,2,4-Trimethylbenzene         | 0.00363            | 0.00292         | 5.4     |
|                  | 1,2-Dibromo-3-chloropropane    | 0.000735 U         | 0.000683 U      | NC      |
|                  | 1,2-Dibromoethane (EDB)        | 0.000299 U         | 0.000278 U      | NC      |
|                  | 1,2-Dichlorobenzene            | 0.000555 U         | 0.000516 U      | NC      |
|                  | 1,2-Dichloroethane (EDC)       | 0.000670 U         | 0.000623 U      | NC      |
|                  | 1,2-Dichloropropane            | 0.000534 U         | 0.000497 U      | NC      |
|                  | 1,3,5-Trimethylbenzene         | 0.0022 J           | 0.00172 J       | 6.1     |
|                  | 1,3-Dichlorobenzene            | 0.000424 U         | 0.000394 U      | NC      |
|                  | 1,3-Dichloropropane            | 0.000390 U         | 0.000362 U      | NC      |
|                  | 1,4-Dichlorobenzene            | 0.000393 U         | 0.000365 U      | NC      |
|                  | 1-Methylnaphthalene            | 0.000982 U         | 0.000913 U      | NC      |
|                  | 2,2-Dichloropropane            | 0.000434 U         | 0.000404 U      | NC      |
|                  | 2-Butanone                     | 0.000801 U         | 0.000745 U      | NC      |
|                  | 2-Chlorotoluene                | 0.000488 U         | 0.000454 U      | NC      |
|                  | 2-Hexanone                     | 0.000520 U         | 0.000484 U      | NC      |
|                  | 2-Methylnaphthalene            | 0.000977 U         | 0.000909 U      | NC      |
|                  | 4-Chlorotoluene                | 0.000482 U         | 0.000448 U      | NC      |
|                  | 4-Isopropyltoluene             | 0.000971 U         | 0.000902 U      | NC      |
|                  | 4-Methyl-2-pentanone           | 0.001056 U         | 0.000982 U      | NC      |
|                  | Acetone                        | 0.00589 J          | 0.00256 J       | 19.7    |
|                  | Benzene                        | 0.0226             | 0.0206          | 2.3     |
|                  | Bromobenzene                   | 0.000696 U         | 0.000647 U      | NC      |
|                  | Bromodichloromethane           | 0.000338 U         | 0.000314 U      | NC      |
|                  | Bromoform                      | 0.000648 U         | 0.000603 U      | NC      |
|                  | Bromomethane                   | 0.001479 U         | 0.001375 U      | NC      |
|                  | Carbon disulfide               | 0.001132 U         | 0.001052 U      | NC      |
|                  | Carbon tetrachloride           | 0.000453 U         | 0.000421 U      | NC      |
|                  | Chlorobenzene                  | 0.000254 U         | 0.000236 U      | NC      |
|                  | Chloroethane                   | 0.002501 U         | 0.002325 U      | NC      |
|                  | Chloroform                     | 0.000515 U         | 0.000479 U      | NC      |
|                  | Chloromethane                  | 0.000295 U         | 0.000274 U      | NC      |
|                  | cis-1,2-DCE                    | 0.000661 U         | 0.000615 U      | NC      |
|                  | cis-1,3-Dichloropropene        | 0.002501 U         | 0.002325 U      | NC      |
|                  | Dibromochloromethane           | 0.000281 U         | 0.000261 U      | NC      |
|                  | Dibromomethane                 | 0.001419 U         | 0.001320 U      | NC      |
|                  | Dichlorodifluoromethane        | 0.001021 U         | 0.000949 U      | NC      |
|                  | Ethylbenzene                   | 0.0035             | 0.00328         | 1.6     |
|                  | Hexachlorobutadiene            | 0.000505 U         | 0.000469 U      | NC      |
|                  | Isopropylbenzene               | 0.000580 U         | 0.000539 U      | NC      |
|                  | Methyl tert-butyl ether (MTBE) | 0.001049 U         | 0.000975 U      | NC      |
|                  | Methylene chloride             | 0.000618 U         | 0.000721 J      | NC      |
|                  | Naphthalene                    | 0.00103 J          | 0.000837 J      | 5.2     |

Table A-3  
Field Duplicate Summary  
SWMU 10 Investigation Report  
Western Refining Southwest, Inc. - Gallup Refinery

|                    | Parameter                   | SWMU 10-6 (10-12') | SWMU 10 DUP03   | RPD (%) |
|--------------------|-----------------------------|--------------------|-----------------|---------|
|                    |                             | Sample Result      | Field Duplicate |         |
|                    | n-Butylbenzene              | 0.001027 U         | 0.000955 U      | NC      |
|                    | n-Propylbenzene             | 0.000638 J         | 0.000477 J      | 7.2     |
|                    | sec-Butylbenzene            | 0.000925 J         | 0.000849 J      | 2.1     |
|                    | Styrene                     | 0.000414 U         | 0.000385 U      | NC      |
|                    | tert-Butylbenzene           | 0.000383 U         | 0.000356 U      | NC      |
|                    | Tetrachloroethene (PCE)     | 0.000496 U         | 0.000461 U      | NC      |
|                    | Toluene                     | 0.0146             | 0.0139          | 1.2     |
|                    | trans-1,2-DCE               | 0.000404 U         | 0.000376 U      | NC      |
|                    | trans-1,3-Dichloropropene   | 0.000327 U         | 0.000304 U      | NC      |
|                    | Trichloroethene (TCE)       | 0.000497 U         | 0.000462 U      | NC      |
|                    | Trichlorofluoromethane      | 0.000505 U         | 0.000470 U      | NC      |
|                    | Vinyl chloride              | 0.000286 U         | 0.000266 U      | NC      |
|                    | Xylenes, Total              | 0.0116             | 0.00984         | 4.1     |
| SVOCs (mg/kg-dry): | 1,2,4-Trichlorobenzene      | 0.120870 U         | 0.112566 U      | NC      |
|                    | 1,2-Dichlorobenzene         | 0.112940 U         | 0.105182 U      | NC      |
|                    | 1,3-Dichlorobenzene         | 0.104874 U         | 0.097669 U      | NC      |
|                    | 1,4-Dichlorobenzene         | 0.123504 U         | 0.115020 U      | NC      |
|                    | 1-Methylnaphthalene         | 0.109448 U         | 0.101929 U      | NC      |
|                    | 2,4,5-Trichlorophenol       | 0.137603 U         | 0.128149 U      | NC      |
|                    | 2,4,6-Trichlorophenol       | 0.136511 U         | 0.127132 U      | NC      |
|                    | 2,4-Dichlorophenol          | 0.118076 U         | 0.109965 U      | NC      |
|                    | 2,4-Dimethylphenol          | 0.092814 U         | 0.086437 U      | NC      |
|                    | 2,4-Dinitrophenol           | 0.049062 U         | 0.045692 U      | NC      |
|                    | 2,4-Dinitrotoluene          | 0.105635 U         | 0.098378 U      | NC      |
|                    | 2,6-Dinitrotoluene          | 0.131249 U         | 0.122232 U      | NC      |
|                    | 2-Chloronaphthalene         | 0.126505 U         | 0.117815 U      | NC      |
|                    | 2-Chlorophenol              | 0.110191 U         | 0.102621 U      | NC      |
|                    | 2-Methylnaphthalene         | 0.108055 U         | 0.100632 U      | NC      |
|                    | 2-Methylphenol              | 0.118374 U         | 0.110242 U      | NC      |
|                    | 2-Nitroaniline              | 0.134357 U         | 0.125127 U      | NC      |
|                    | 2-Nitrophenol               | 0.106127 U         | 0.098836 U      | NC      |
|                    | 3,3'-Dichlorobenzidine      | 0.090922 U         | 0.084676 U      | NC      |
|                    | 3+4-Methylphenol            | 0.122378 U         | 0.113970 U      | NC      |
|                    | 3-Nitroaniline              | 0.113765 U         | 0.105949 U      | NC      |
|                    | 4,6-Dinitro-2-methylphenol  | 0.064203 U         | 0.059792 U      | NC      |
|                    | 4-Bromophenyl phenyl ether  | 0.125276 U         | 0.116669 U      | NC      |
|                    | 4-Chloro-3-methylphenol     | 0.117566 U         | 0.109489 U      | NC      |
|                    | 4-Chloroaniline             | 0.109198 U         | 0.101696 U      | NC      |
|                    | 4-Chlorophenyl phenyl ether | 0.175825 U         | 0.163746 U      | NC      |
|                    | 4-Nitroaniline              | 0.106056 U         | 0.098770 U      | NC      |
|                    | 4-Nitrophenol               | 0.101733 U         | 0.094744 U      | NC      |
|                    | Acenaphthene                | 0.143884 U         | 0.133999 U      | NC      |
|                    | Acenaphthylene              | 0.119613 U         | 0.111396 U      | NC      |
|                    | Aniline                     | 0.102777 U         | 0.095717 U      | NC      |
|                    | Anthracene                  | 0.100297 U         | 0.093407 U      | NC      |
|                    | Azobenzene                  | 0.132366 U         | 0.123272 U      | NC      |
|                    | Benz(a)anthracene           | 0.092803 U         | 0.086427 U      | NC      |
|                    | Benzo(a)pyrene              | 0.132513 U         | 0.123409 U      | NC      |
|                    | Benzo(b)fluoranthene        | 0.112124 U         | 0.104421 U      | NC      |
|                    | Benzo(g,h,i)perylene        | 0.134367 U         | 0.125136 U      | NC      |
|                    | Benzo(k)fluoranthene        | 0.133438 U         | 0.124271 U      | NC      |
|                    | Benzoic acid                | 0.067280 U         | 0.062658 U      | NC      |
|                    | Benzyl alcohol              | 0.104881 U         | 0.097676 U      | NC      |
|                    | Bis(2-chloroethoxy)methane  | 0.112695 U         | 0.104953 U      | NC      |
|                    | Bis(2-chloroethyl)ether     | 0.116670 U         | 0.108655 U      | NC      |

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                     | Parameter                   | SWMU 10-6 (10-12') | SWMU 10 DUP03   | RPD (%) |
|---------------------|-----------------------------|--------------------|-----------------|---------|
|                     |                             | Sample Result      | Field Duplicate |         |
|                     | Bis(2-chloroisopropyl)ether | 0.090206 U         | 0.084009 U      | NC      |
|                     | Bis(2-ethylhexyl)phthalate  | 0.132656 U         | 0.123542 U      | NC      |
|                     | Butyl benzyl phthalate      | 0.144149 U         | 0.134246 U      | NC      |
|                     | Carbazole                   | 0.109971 U         | 0.102416 U      | NC      |
|                     | Chrysene                    | 0.118564 U         | 0.110419 U      | NC      |
|                     | Dibenz(a,h)anthracene       | 0.127923 U         | 0.119135 U      | NC      |
|                     | Dibenzofuran                | 0.126280 U         | 0.117604 U      | NC      |
|                     | Diethyl phthalate           | 0.134003 U         | 0.124797 U      | NC      |
|                     | Dimethyl phthalate          | 0.108666 U         | 0.101200 U      | NC      |
|                     | Di-n-butyl phthalate        | 0.133536 U         | 0.124362 U      | NC      |
|                     | Di-n-octyl phthalate        | 0.131446 U         | 0.122416 U      | NC      |
|                     | Fluoranthene                | 0.143283 U         | 0.133440 U      | NC      |
|                     | Fluorene                    | 0.160744 U         | 0.149701 U      | NC      |
|                     | Hexachlorobenzene           | 0.111443 U         | 0.103787 U      | NC      |
|                     | Hexachlorobutadiene         | 0.116817 U         | 0.108792 U      | NC      |
|                     | Hexachlorocyclopentadiene   | 0.081898 U         | 0.076272 U      | NC      |
|                     | Hexachloroethane            | 0.105103 U         | 0.097883 U      | NC      |
|                     | Indeno(1,2,3-cd)pyrene      | 0.134185 U         | 0.124967 U      | NC      |
|                     | Isophorone                  | 0.125890 U         | 0.117241 U      | NC      |
|                     | Naphthalene                 | 0.112350 U         | 0.104631 U      | NC      |
|                     | Nitrobenzene                | 0.121569 U         | 0.113217 U      | NC      |
|                     | N-Nitrosodi-n-propylamine   | 0.122390 U         | 0.113982 U      | NC      |
|                     | N-Nitrosodiphenylamine      | 0.105712 U         | 0.098449 U      | NC      |
|                     | Pentachlorophenol           | 0.073720 U         | 0.068656 U      | NC      |
|                     | Phenanthrene                | 0.123526 U         | 0.115040 U      | NC      |
|                     | Phenol                      | 0.105819 U         | 0.098549 U      | NC      |
|                     | Pyrene                      | 0.152390 U         | 0.141921 U      | NC      |
|                     | Pyridine                    | 0.099219 U         | 0.092403 U      | NC      |
| Metals (mg/kg-dry): | Antimony                    | 1.9 J              | 1.9 J           | 0.0     |
|                     | Arsenic                     | 2.1 J              | 2.2 J           | 1.2     |
|                     | Barium                      | 550                | 450             | 5.0     |
|                     | Beryllium                   | 1                  | 0.95            | 1.3     |
|                     | Cadmium                     | 0.033 U            | 0.03 U          | NC      |
|                     | Chromium                    | 14                 | 13              | 1.9     |
|                     | Cobalt                      | 8                  | 7               | 3.3     |
|                     | Cyanide                     | 0.296 U            | 0.28 U          | NC      |
|                     | Hexavalent Chromium         | 2,368 U            | 260             | NC      |
|                     | Iron                        | 20000              | 20000           | 0.0     |
|                     | Lead                        | 1.6                | 1.4             | 3.3     |
|                     | Manganese                   | 760                | 730             | 1.0     |
|                     | Mercury                     | 0.0035 U           | 0.0031 U        | NC      |
|                     | Nickel                      | 16                 | 13              | 5.2     |
|                     | Selenium                    | 1.821 U            | 1.69 U          | NC      |
|                     | Silver                      | 0.035 U            | 0.03 U          | NC      |
|                     | Vanadium                    | 14                 | 17              | 4.8     |
|                     | Zinc                        | 23                 | 21              | 2.3     |

**Notes:**

RPD = Relative percent difference; [(difference)/(average)]\* 100  
NC = Not calculated; RPD values were not calculated for non-detects  
ug/kg-dry = micrograms per kilogram dry  
mg/kg-dry = milligrams per kilogram  
bold value = Field Duplicate RPD Outlier

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                  | Parameter                      | SWMU 10-17 (6-8')<br>Sample Result | SWMU 10 DUP04<br>Field Duplicate | RPD<br>(%) |
|------------------|--------------------------------|------------------------------------|----------------------------------|------------|
| TPH (mg/kg-dry): | Gasoline Range Organics (GRO)  | 1300                               | 1100 v                           | 4.2        |
|                  | Diesel Range Organics (DRO)    | 7800                               | 3000 v                           | 22.2       |
|                  | Motor Oil Range Organics (MRO) | ND U                               | < 488 u                          | NC         |
| VOCs (ug/kg-dry) | 1,1,1,2-Tetrachloroethane      | ND U                               | < 0.0526 u                       | NC         |
|                  | 1,1,1-Trichloroethane          | ND U                               | < 0.0975 u                       | NC         |
|                  | 1,1,2,2-Tetrachloroethane      | ND U                               | < 0.0778 u                       | NC         |
|                  | 1,1,2-Trichloroethane          | ND U                               | < 0.0771 u                       | NC         |
|                  | 1,1-Dichloroethane             | ND U                               | < 0.2568 u                       | NC         |
|                  | 1,1-Dichloroethene             | ND U                               | < 0.0442 u                       | NC         |
|                  | 1,1-Dichloropropene            | ND U                               | < 0.0956 u                       | NC         |
|                  | 1,2,3-Trichlorobenzene         | ND U                               | < 0.137 u                        | NC         |
|                  | 1,2,3-Trichloropropane         | ND U                               | < 0.1077 u                       | NC         |
|                  | 1,2,4-Trichlorobenzene         | ND U                               | < 0.11844 u                      | NC         |
|                  | 1,2,4-Trimethylbenzene         | 27                                 | 17 v                             | 11.4       |
|                  | 1,2-Dibromo-3-chloropropane    | ND U                               | < 0.1103 u                       | NC         |
|                  | 1,2-Dibromoethane (EDB)        | ND U                               | < 0.0511 u                       | NC         |
|                  | 1,2-Dichlorobenzene            | ND U                               | < 0.03871 u                      | NC         |
|                  | 1,2-Dichloroethane (EDC)       | ND U                               | < 0.2568 u                       | NC         |
|                  | 1,2-Dichloropropane            | ND U                               | < 0.0359 u                       | NC         |
|                  | 1,3,5-Trimethylbenzene         | 10                                 | 6.2 v                            | 11.7       |
|                  | 1,3-Dichlorobenzene            | ND U                               | < 0.07558 u                      | NC         |
|                  | 1,3-Dichloropropane            | ND U                               | < 0.0884 u                       | NC         |
|                  | 1,4-Dichlorobenzene            | ND U                               | < 0.07995 u                      | NC         |
|                  | 1-Methylnaphthalene            | 48                                 | 31 v                             | 10.8       |
|                  | 2,2-Dichloropropane            | ND U                               | < 0.0934 u                       | NC         |
|                  | 2-Butanone                     | ND U                               | < 0.4127 u                       | NC         |
|                  | 2-Chlorotoluene                | ND U                               | < 0.0725 u                       | NC         |
|                  | 2-Hexanone                     | ND U                               | < 0.1579 u                       | NC         |
|                  | 2-Methylnaphthalene            | 87                                 | 60 v                             | 9.2        |
|                  | 4-Chlorotoluene                | ND U                               | < 0.0485 u                       | NC         |
|                  | 4-Isopropyltoluene             | 1.6                                | 1.1 v                            | 9.3        |
|                  | 4-Methyl-2-pentanone           | ND U                               | < 0.1114 u                       | NC         |
|                  | Acetone                        | ND U                               | < 0.4786 u                       | NC         |
|                  | Benzene                        | 1.9                                | 1.2 v                            | 11.3       |
|                  | Bromobenzene                   | ND U                               | < 0.0476 u                       | NC         |
|                  | Bromodichloromethane           | ND U                               | < 0.0525 u                       | NC         |
|                  | Bromoform                      | ND U                               | < 0.0516 u                       | NC         |
|                  | Bromomethane                   | ND U                               | < 0.1239 u                       | NC         |
|                  | Carbon disulfide               | ND U                               | < 0.4619 u                       | NC         |
|                  | Carbon tetrachloride           | ND U                               | < 0.0625 u                       | NC         |
|                  | Chlorobenzene                  | ND U                               | < 0.0494 u                       | NC         |
|                  | Chloroethane                   | ND U                               | < 0.4969 u                       | NC         |
|                  | Chloroform                     | ND U                               | < 0.0669 u                       | NC         |
|                  | Chloromethane                  | ND U                               | < 0.0743 u                       | NC         |
|                  | cis-1,2-DCE                    | ND U                               | < 0.0812 u                       | NC         |
|                  | cis-1,3-Dichloropropene        | ND U                               | < 0.0445 u                       | NC         |
|                  | Dibromochloromethane           | ND U                               | < 0.0481 u                       | NC         |
|                  | Dibromomethane                 | ND U                               | < 0.0676 u                       | NC         |
|                  | Dichlorodifluoromethane        | ND U                               | < 0.1614 u                       | NC         |
|                  | Ethylbenzene                   | 11                                 | 8.3 v                            | 7.0        |
|                  | Hexachlorobutadiene            | ND U                               | < 0.08913 u                      | NC         |
|                  | Isopropylbenzene               | 3.1                                | 2.3 v                            | 7.4        |
|                  | Methyl tert-butyl ether (MTBE) | ND U                               | < 0.0908 u                       | NC         |
|                  | Methylene chloride             | ND U                               | < 0.2568 u                       | NC         |

Table A-3  
Field Duplicate Summary  
SWMU 10 Investigation Report  
Western Refining Southwest, Inc. - Gallup Refinery

|                    | Parameter                   | SWMU 10-17 (6-8')<br>Sample Result | SWMU 10 DUP04<br>Field Duplicate | RPD<br>(%) |
|--------------------|-----------------------------|------------------------------------|----------------------------------|------------|
|                    | Naphthalene                 | 20                                 | 13 v                             | 10.6       |
|                    | n-Butylbenzene              | 2.8                                | 1.9 J                            | 9.6        |
|                    | n-Propylbenzene             | 5.2                                | 3.3 v                            | 11.2       |
|                    | sec-Butylbenzene            | 1.8                                | 1 v                              | 14.3       |
|                    | Styrene                     | ND                                 | U < 0.1178                       | u NC       |
|                    | tert-Butylbenzene           | ND                                 | U < 0.0735                       | u NC       |
|                    | Tetrachloroethene (PCE)     | ND                                 | U < 0.0556                       | u NC       |
|                    | Toluene                     | 34                                 | 23 v                             | 9.6        |
|                    | trans-1,2-DCE               | ND                                 | U < 0.0417                       | u NC       |
|                    | trans-1,3-Dichloropropene   | ND                                 | U < 0.1156                       | u NC       |
|                    | Trichloroethene (TCE)       | ND                                 | U < 0.0625                       | u NC       |
|                    | Trichlorofluoromethane      | ND                                 | U < 0.1938                       | u NC       |
|                    | Vinyl chloride              | ND                                 | U < 0.2568                       | u NC       |
|                    | Xylenes, Total              | 86                                 | 65 v                             | 7.0        |
| SVOCs (mg/kg-dry): | 1,2,4-Trichlorobenzene      | ND                                 | U < 1.163                        | u NC       |
|                    | 1,2-Dichlorobenzene         | ND                                 | U < 1.0867                       | u NC       |
|                    | 1,3-Dichlorobenzene         | ND                                 | U < 1.0091                       | u NC       |
|                    | 1,4-Dichlorobenzene         | ND                                 | U < 1.1883                       | u NC       |
|                    | 1-Methylnaphthalene         | 76                                 | 28 v                             | 23.1       |
|                    | 2,4,5-Trichlorophenol       | ND                                 | U < 1.324                        | u NC       |
|                    | 2,4,6-Trichlorophenol       | ND                                 | U < 1.3135                       | u NC       |
|                    | 2,4-Dichlorophenol          | ND                                 | U < 1.1361                       | u NC       |
|                    | 2,4-Dimethylphenol          | 36                                 | 18 v                             | 16.7       |
|                    | 2,4-Dinitrophenol           | ND                                 | U < 0.4721                       | u NC       |
|                    | 2,4-Dinitrotoluene          | ND                                 | U < 1.0164                       | u NC       |
|                    | 2,6-Dinitrotoluene          | ND                                 | U < 1.2629                       | u NC       |
|                    | 2-Chloronaphthalene         | ND                                 | U < 1.2172                       | u NC       |
|                    | 2-Chlorophenol              | ND                                 | U < 1.0602                       | u NC       |
|                    | 2-Methylnaphthalene         | 130                                | 45 v                             | 24.3       |
|                    | 2-Methylphenol              | 56                                 | 29 v                             | 15.9       |
|                    | 2-Nitroaniline              | ND                                 | U < 1.2928                       | u NC       |
|                    | 2-Nitrophenol               | ND                                 | U < 1.0211                       | u NC       |
|                    | 3,3'-Dichlorobenzidine      | ND                                 | U < 0.8748                       | u NC       |
|                    | 3+4-Methylphenol            | 100                                | 60 v                             | 12.5       |
|                    | 3-Nitroaniline              | ND                                 | U < 1.0946                       | u NC       |
|                    | 4,6-Dinitro-2-methylphenol  | ND                                 | U < 0.6178                       | u NC       |
|                    | 4-Bromophenyl phenyl ether  | ND                                 | U < 1.2054                       | u NC       |
|                    | 4-Chloro-3-methylphenol     | ND                                 | U < 1.1312                       | u NC       |
|                    | 4-Chloroaniline             | ND                                 | U < 1.0507                       | u NC       |
|                    | 4-Chlorophenyl phenyl ether | ND                                 | U < 1.6918                       | u NC       |
|                    | 4-Nitroaniline              | ND                                 | U < 1.0205                       | u NC       |
|                    | 4-Nitrophenol               | ND                                 | U < 0.9789                       | u NC       |
|                    | Acenaphthene                | ND                                 | U < 1.3844                       | u NC       |
|                    | Acenaphthylene              | ND                                 | U < 1.1509                       | u NC       |
|                    | Aniline                     | ND                                 | U < 0.9889                       | u NC       |
|                    | Anthracene                  | ND                                 | U < 0.965                        | u NC       |
|                    | Azobenzene                  | ND                                 | U < 1.2736                       | u NC       |
|                    | Benz(a)anthracene           | ND                                 | U < 0.8929                       | u NC       |
|                    | Benzo(a)pyrene              | ND                                 | U < 1.275                        | u NC       |
|                    | Benzo(b)fluoranthene        | ND                                 | U < 1.0788                       | u NC       |
|                    | Benzo(g,h,i)perylene        | ND                                 | U < 1.2929                       | u NC       |
|                    | Benzo(k)fluoranthene        | ND                                 | U < 1.2839                       | u NC       |
|                    | Benzoic acid                | ND                                 | U < 0.6474                       | u NC       |
|                    | Benzyl alcohol              | ND                                 | U < 1.0092                       | u NC       |

Table A-3  
Field Duplicate Summary  
SWMU 10 Investigation Report  
Western Refining Southwest, Inc. - Gallup Refinery

|                     | Parameter                   | SWMU 10-17 (6-8')<br>Sample Result | SWMU 10 DUP04<br>Field Duplicate | RPD<br>(%) |
|---------------------|-----------------------------|------------------------------------|----------------------------------|------------|
|                     | Bis(2-chloroethoxy)methane  | ND U                               | < 1.0843 u                       | NC         |
|                     | Bis(2-chloroethyl)ether     | ND U                               | < 1.1226 u                       | NC         |
|                     | Bis(2-chloroisopropyl)ether | ND U                               | < 0.8679 u                       | NC         |
|                     | Bis(2-ethylhexyl)phthalate  | ND U                               | < 1.2764 u                       | NC         |
|                     | Butyl benzyl phthalate      | ND U                               | < 1.387 u                        | NC         |
|                     | Carbazole                   | ND U                               | < 1.0581 u                       | NC         |
|                     | Chrysene                    | ND U                               | < 1.1408 u                       | NC         |
|                     | Dibenz(a,h)anthracene       | ND U                               | < 1.2309 u                       | NC         |
|                     | Dibenzofuran                | ND U                               | < 1.215 u                        | NC         |
|                     | Diethyl phthalate           | ND U                               | < 1.2894 u                       | NC         |
|                     | Dimethyl phthalate          | ND U                               | < 1.0456 u                       | NC         |
|                     | Di-n-butyl phthalate        | ND U                               | < 1.2849 u                       | NC         |
|                     | Di-n-octyl phthalate        | ND U                               | < 1.2648 u                       | NC         |
|                     | Fluoranthene                | ND U                               | < 1.3787 u                       | NC         |
|                     | Fluorene                    | 7.1                                | 3.2 v                            | 18.9       |
|                     | Hexachlorobenzene           | ND U                               | < 1.0723 u                       | NC         |
|                     | Hexachlorobutadiene         | ND U                               | < 1.124 u                        | NC         |
|                     | Hexachlorocyclopentadiene   | ND U                               | < 0.788 u                        | NC         |
|                     | Hexachloroethane            | ND U                               | < 1.0113 u                       | NC         |
|                     | Indeno(1,2,3-cd)pyrene      | ND U                               | < 1.2911 u                       | NC         |
|                     | Naphthalene                 | 24                                 | 8.8 v                            | 23.2       |
|                     | Nitrobenzene                | ND U                               | < 1.1697 u                       | NC         |
|                     | N-Nitrosodi-n-propylamine   | ND U                               | < 1.1776 u                       | NC         |
|                     | N-Nitrosodiphenylamine      | ND U                               | < 1.0171 u                       | NC         |
|                     | Pentachlorophenol           | ND U                               | < 0.7093 u                       | NC         |
|                     | Phenanthrene                | 15                                 | 6 v                              | 21.4       |
|                     | Phenol                      | 52                                 | 26 v                             | 16.7       |
|                     | Pyrene                      | ND U                               | < 1.4663 u                       | NC         |
|                     | Pyridine                    | ND U                               | < 0.9547 u                       | NC         |
| Metals (mg/kg-dry): | Antimony                    | ND U                               | < 1.5264 u                       | NC         |
|                     | Arsenic                     | 1.5 J                              | 1.1 J                            | 7.7        |
|                     | Barium                      | 200                                | 160 v                            | 5.6        |
|                     | Beryllium                   | 1                                  | 1.1 v                            | 2.4        |
|                     | Cadmium                     | ND U                               | < 0.0322 u                       | NC         |
|                     | Chromium                    | 11                                 | 11 v                             | 0.0        |
|                     | Cobalt                      | 5.3                                | < 2.2887 u                       | NC         |
|                     | Cyanide                     | ND U                               | 5.4 v                            | NC         |
|                     | Hexavalent Chromium         | ND U                               | < 0.2861 u                       | NC         |
|                     | Iron                        | 16000                              | 16000 v                          | 0.0        |
|                     | Lead                        | 4.4                                | 5.3 v                            | 4.6        |
|                     | Manganese                   | 210                                | 200 v                            | 1.2        |
|                     | Mercury                     | ND U                               | < 0.0032 u                       | NC         |
|                     | Nickel                      | 9.6                                | 9.5 v                            | 0.3        |
|                     | Selenium                    | ND U                               | < 1.7774 u                       | NC         |
|                     | Silver                      | ND U                               | < 0.0345 u                       | NC         |
|                     | Vanadium                    | 19                                 | 19 v                             | 0.0        |
|                     | Zinc                        | 17                                 | 17 v                             | 0.0        |

**Notes:**

RPD = Relative percent difference; [(difference)/(average)]\* 100

NC = Not calculated; RPD values were not calculated for non-detects

ug/kg-dry = micrograms per kilogram dry

mg/kg-dry = milligrams per kilogram

bold value = Field Duplicate RPD Outlier

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                  | Parameter                      | SWMU 10-22 (2-2.5') |   | DUP01           |   |      |
|------------------|--------------------------------|---------------------|---|-----------------|---|------|
|                  |                                | 1605943-004         |   | 1605943-006     |   | RPD  |
|                  |                                | Sample Result       |   | Field Duplicate |   | %    |
| TPH (mg/kg-dry)  | Gasoline Range Organics (GRO)  | 0.6                 | J | < 0.858         | u | NC   |
|                  | Diesel Range Organics (DRO)    | < 4.5               | u | 150             | v | NC   |
|                  | Motor Oil Range Organics (MRO) | < 62                | u | 310             | v | NC   |
| VOCs (mg/kg-dry) | 1,1,1,2-Tetrachloroethane      | < 0.0017            | u | < 0.0003        | u | NC   |
|                  | 1,1,1-Trichloroethane          | < 0.0017            | u | < 0.0002        | u | NC   |
|                  | 1,1,2,2-Tetrachloroethane      | < 0.0017            | u | < 0.0004        | u | NC   |
|                  | 1,1,2-Trichloroethane          | < 0.0017            | u | < 0.0016        | u | NC   |
|                  | 1,1-Dichloroethane             | < 0.0017            | u | < 0.0004        | u | NC   |
|                  | 1,1-Dichloroethene             | < 0.0003            | u | < 0.0003        | u | NC   |
|                  | 1,1-Dichloropropene            | < 0.0017            | u | < 0.0003        | u | NC   |
|                  | 1,2,3-Trichlorobenzene         | < 0.0004            | u | < 0.0006        | u | NC   |
|                  | 1,2,3-Trichloropropane         | < 0.0017            | u | < 0.0006        | u | NC   |
|                  | 1,2,4-Trichlorobenzene         | < 0.0005            | u | < 0.00059       | u | NC   |
|                  | 1,2,4-Trimethylbenzene         | 0.0003              | J | 0.000972        | J | 26.4 |
|                  | 1,2-Dibromo-3-chloropropane    | < 0.0002            | u | < 0.0005        | u | NC   |
|                  | 1,2-Dibromoethane (EDB)        | < 0.0017            | u | < 0.0002        | u | NC   |
|                  | 1,2-Dichlorobenzene            | < 0.0002            | u | < 0.00035       | u | NC   |
|                  | 1,2-Dichloroethane (EDC)       | < 0.0017            | u | < 0.0004        | u | NC   |
|                  | 1,2-Dichloropropane            | < 0.0017            | u | < 0.0003        | u | NC   |
|                  | 1,3,5-Trimethylbenzene         | < 0.0003            | u | 0.000422        | J | NC   |
|                  | 1,3-Dichlorobenzene            | < 0.0003            | u | < 0.00027       | u | NC   |
|                  | 1,3-Dichloropropane            | < 0.0017            | u | < 0.0002        | u | NC   |
|                  | 1,4-Dichlorobenzene            | < 0.0003            | u | < 0.00025       | u | NC   |
|                  | 1-Methylnaphthalene            | < 0.0002            | u | < 0.00062       | u | NC   |
|                  | 2,2-Dichloropropane            | < 0.0002            | u | < 0.0003        | u | NC   |
|                  | 2-Butanone                     | < 0.0006            | u | 0.00253         | J | NC   |
|                  | 2-Chlorotoluene                | < 0.0003            | u | < 0.0003        | u | NC   |
|                  | 2-Hexanone                     | < 0.0004            | u | 0.000653        | J | NC   |
|                  | 2-Methylnaphthalene            | < 0.0004            | u | < 0.00062       | u | NC   |
|                  | 4-Chlorotoluene                | < 0.0003            | u | < 0.0003        | u | NC   |
|                  | 4-Isopropyltoluene             | < 0.0003            | u | < 0.0006        | u | NC   |
|                  | 4-Methyl-2-pentanone           | < 0.0033            | u | < 0.0007        | u | NC   |
|                  | Acetone                        | 0.0117              | v | 0.0166          | v | 8.7  |
|                  | Benzene                        | < 0.0017            | u | 0.00104         | J | NC   |
|                  | Bromobenzene                   | < 0.0002            | u | < 0.0004        | u | NC   |
|                  | Bromodichloromethane           | < 0.0017            | u | < 0.0002        | u | NC   |
|                  | Bromoform                      | < 0.0017            | u | < 0.0004        | u | NC   |
|                  | Bromomethane                   | < 0.0003            | u | < 0.0009        | u | NC   |
|                  | Carbon disulfide               | < 0.0006            | u | < 0.0007        | u | NC   |
|                  | Carbon tetrachloride           | < 0.0017            | u | < 0.0003        | u | NC   |
|                  | Chlorobenzene                  | < 0.0002            | u | < 0.0002        | u | NC   |
|                  | Chloroethane                   | < 0.0003            | u | < 0.0016        | u | NC   |
|                  | Chloroform                     | < 0.0017            | u | < 0.0003        | u | NC   |
|                  | Chloromethane                  | < 0.0004            | u | < 0.0002        | u | NC   |
|                  | cis-1,2-DCE                    | < 0.0017            | u | < 0.0004        | u | NC   |
|                  | cis-1,3-Dichloropropene        | < 0.0017            | u | < 0.0016        | u | NC   |
|                  | Dibromochloromethane           | < 0.0017            | u | < 0.0002        | u | NC   |
|                  | Dibromomethane                 | < 0.0017            | u | < 0.0009        | u | NC   |
|                  | Dichlorodifluoromethane        | < 0.001             | u | < 0.0007        | u | NC   |
|                  | Ethylbenzene                   | < 0.0002            | u | 0.000638        | J | NC   |



**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                   | Parameter                      | SWMU 10-22 (2-2.5') |   | DUP01           |   |      |
|-------------------|--------------------------------|---------------------|---|-----------------|---|------|
|                   |                                | 1605943-004         |   | 1605943-006     |   | RPD  |
|                   |                                | Sample Result       |   | Field Duplicate |   | %    |
|                   | Hexachlorobutadiene            | < 0.0004            | u | < 0.00032       | u | NC   |
|                   | Isopropylbenzene               | < 0.0002            | u | < 0.0004        | u | NC   |
|                   | Methyl tert-butyl ether (MTBE) | < 0.0003            | u | < 0.0007        | u | NC   |
|                   | Methylene chloride             | < 0.0017            | u | 0.00051         | J | NC   |
|                   | Naphthalene                    | < 0.0017            | u | < 0.00052       | u | NC   |
|                   | n-Butylbenzene                 | < 0.0004            | u | < 0.0007        | u | NC   |
|                   | n-Propylbenzene                | < 0.0003            | u | < 0.0003        | u | NC   |
|                   | sec-Butylbenzene               | < 0.0003            | u | 0.000534        | J | NC   |
|                   | Styrene                        | < 0.0002            | u | < 0.0003        | u | NC   |
|                   | tert-Butylbenzene              | < 0.0002            | u | < 0.0002        | u | NC   |
|                   | Tetrachloroethene (PCE)        | < 0.0002            | u | < 0.0003        | u | NC   |
|                   | Toluene                        | 0.0012              | J | 0.00232         | v | 15.9 |
|                   | trans-1,2-DCE                  | < 0.0002            | u | < 0.0003        | u | NC   |
|                   | trans-1,3-Dichloropropene      | < 0.0002            | u | < 0.0002        | u | NC   |
|                   | Trichloroethene (TCE)          | < 0.0017            | u | < 0.0003        | u | NC   |
|                   | Trichlorofluoromethane         | < 0.0002            | u | < 0.0003        | u | NC   |
|                   | Vinyl chloride                 | < 0.0004            | u | < 0.0002        | u | NC   |
|                   | Xylenes, Total                 | < 0.0007            | u | 0.00171         | v | NC   |
| SVOCs (mg/kg-dry) | 1,2,4-Trichlorobenzene         | < 0.13              | u | < 2.8155        | u | NC   |
|                   | 1,2-Dichlorobenzene            | < 0.094             | u | < 2.6308        | u | NC   |
|                   | 1,3-Dichlorobenzene            | < 0.095             | u | < 2.4429        | u | NC   |
|                   | 1,4-Dichlorobenzene            | < 0.1               | u | < 2.8768        | u | NC   |
|                   | 1-Methylnaphthalene            | < 0.12              | u | < 2.5494        | u | NC   |
|                   | 2,4,5-Trichlorophenol          | < 0.12              | u | < 3.2052        | u | NC   |
|                   | 2,4,6-Trichlorophenol          | < 0.1               | u | < 3.1798        | u | NC   |
|                   | 2,4-Dichlorophenol             | < 0.11              | u | < 2.7504        | u | NC   |
|                   | 2,4-Dimethylphenol             | < 0.13              | u | < 2.1619        | u | NC   |
|                   | 2,4-Dinitrophenol              | < 0.081             | u | < 1.1428        | u | NC   |
|                   | 2,4-Dinitrotoluene             | < 0.11              | u | < 2.4606        | u | NC   |
|                   | 2,6-Dinitrotoluene             | < 0.13              | u | < 3.0572        | u | NC   |
|                   | 2-Chloronaphthalene            | < 0.097             | u | < 2.9467        | u | NC   |
|                   | 2-Chlorophenol                 | < 0.097             | u | < 2.5667        | u | NC   |
|                   | 2-Methylnaphthalene            | < 0.15              | u | < 2.517         | u | NC   |
|                   | 2-Methylphenol (cresol,o-)     | < 0.1               | u | < 2.7573        | u | NC   |
|                   | 2-Nitroaniline                 | < 0.13              | u | < 3.1296        | u | NC   |
|                   | 2-Nitrophenol                  | < 0.12              | u | < 2.4721        | u | NC   |
|                   | 3,3'-Dichlorobenzidine         | < 0.09              | u | < 2.1179        | u | NC   |
|                   | 3+4-Methylphenol               | < 0.089             | u | < 2.8506        | u | NC   |
|                   | 3-Nitroaniline                 | < 0.11              | u | < 2.65          | u | NC   |
|                   | 4,6-Dinitro-2-methylphenol     | < 0.074             | u | < 1.4955        | u | NC   |
|                   | 4-Bromophenyl phenyl ether     | < 0.12              | u | < 2.9181        | u | NC   |
|                   | 4-Chloro-3-methylphenol        | < 0.15              | u | < 2.7385        | u | NC   |
|                   | 4-Chloroaniline                | < 0.13              | u | < 2.5436        | u | NC   |
|                   | 4-Chlorophenyl phenyl ether    | < 0.14              | u | < 4.0956        | u | NC   |
|                   | 4-Nitroaniline                 | < 0.086             | u | < 2.4704        | u | NC   |
|                   | 4-Nitrophenol                  | < 0.093             | u | < 2.3697        | u | NC   |
|                   | Acenaphthene                   | < 0.11              | u | < 3.3515        | u | NC   |
|                   | Acenaphthylene                 | < 0.1               | u | < 2.7862        | u | NC   |
|                   | Aniline                        | < 0.12              | u | < 2.394         | u | NC   |
|                   | Anthracene                     | < 0.081             | u | < 2.3363        | u | NC   |

Table A-3  
Field Duplicate Summary  
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Western Refining Southwest, Inc. - Gallup Refinery

|                     | Parameter                   | SWMU 10-22 (2-2.5') |       | DUP01           |          |      |     |
|---------------------|-----------------------------|---------------------|-------|-----------------|----------|------|-----|
|                     |                             | 1605943-004         |       | 1605943-006     |          | RPD  |     |
|                     |                             | Sample Result       |       | Field Duplicate |          | %    |     |
|                     | Azobenzene                  | < 0.15              | u     | < 3.0833        | u        | NC   |     |
|                     | Benz(a)anthracene           | < 0.11              | u     | < 2.1617        | u        | NC   |     |
|                     | Benzo(a)pyrene              | < 0.093             | u     | < 3.0867        | u        | NC   |     |
|                     | Benzo(b)fluoranthene        | < 0.11              | u     | < 2.6117        | u        | NC   |     |
|                     | Benzo(g,h,i)perylene        | < 0.11              | u     | < 3.1299        | u        | NC   |     |
|                     | Benzo(k)fluoranthene        | < 0.11              | u     | < 3.1082        | u        | NC   |     |
|                     | Benzoic acid                | < 0.1               | u     | < 1.5672        | u        | NC   |     |
|                     | Benzyl alcohol              | < 0.096             | u     | < 2.443         | u        | NC   |     |
|                     | Bis(2-chloroethoxy)methane  | < 0.13              | u     | < 2.6251        | u        | NC   |     |
|                     | Bis(2-chloroethyl)ether     | < 0.09              | u     | < 2.7176        | u        | NC   |     |
|                     | Bis(2-chloroisopropyl)ether | < 0.11              | u     | < 2.1012        | u        | NC   |     |
|                     | Bis(2-ethylhexyl)phthalate  | 0.19                | J     | < 3.09          | u        | NC   |     |
|                     | Butyl benzyl phthalate      | < 0.11              | u     | < 3.3577        | u        | NC   |     |
|                     | Carbazole                   | < 0.083             | u     | < 2.5616        | u        | NC   |     |
|                     | Chrysene                    | < 0.1               | u     | < 2.7618        | u        | NC   |     |
|                     | Dibenz(a,h)anthracene       | < 0.099             | u     | < 2.9798        | u        | NC   |     |
|                     | Dibenzofuran                | < 0.12              | u     | < 2.9415        | u        | NC   |     |
|                     | Diethyl phthalate           | < 0.12              | u     | < 3.1214        | u        | NC   |     |
|                     | Dimethyl phthalate          | < 0.12              | u     | < 2.5312        | u        | NC   |     |
|                     | Di-n-butyl phthalate        | 0.12                | J     | < 3.1105        | u        | NC   |     |
|                     | Di-n-octyl phthalate        | < 0.1               | u     | < 3.0618        | u        | NC   |     |
|                     | Fluoranthene                | < 0.071             | u     | < 3.3375        | u        | NC   |     |
|                     | Fluorene                    | < 0.11              | u     | < 3.7443        | u        | NC   |     |
|                     | Hexachlorobenzene           | < 0.097             | u     | < 2.5959        | u        | NC   |     |
|                     | Hexachlorobutadiene         | < 0.14              | u     | < 2.7211        | u        | NC   |     |
|                     | Hexachlorocyclopentadiene   | < 0.14              | u     | < 1.9077        | u        | NC   |     |
|                     | Hexachloroethane            | < 0.11              | u     | < 2.4482        | u        | NC   |     |
|                     | Indeno(1,2,3-cd)pyrene      | < 0.096             | u     | < 3.1256        | u        | NC   |     |
|                     | Naphthalene                 | < 0.12              | u     | < 2.617         | u        | NC   |     |
|                     | Nitrobenzene                | < 0.13              | u     | < 2.8318        | u        | NC   |     |
|                     | N-Nitrosodi-n-propylamine   | < 0.12              | u     | < 2.8509        | u        | NC   |     |
|                     | N-Nitrosodiphenylamine      | < 0.12              | u     | < 2.4624        | u        | NC   |     |
|                     | Pentachlorophenol           | < 0.079             | u     | < 1.7172        | u        | NC   |     |
|                     | Phenanthrene                | < 0.083             | u     | < 2.8773        | u        | NC   |     |
|                     | Phenol                      | < 0.092             | u     | < 2.4649        | u        | NC   |     |
|                     | Pyrene                      | < 0.093             | u     | < 3.5497        | u        | NC   |     |
|                     | Pyridine                    | < 0.097             | u     | < 2.3111        | u        | NC   |     |
|                     | Metals (mg/kg-dry)          | Antimony            | < 7.3 | u               | < 1.4517 | u    | NC  |
|                     |                             | Arsenic             | 2.4   | J               | 2.7      | J    | 2.9 |
|                     |                             | Barium              | 290   | v               | 380      | v    | 6.7 |
| Beryllium           |                             | 0.89                | v     | 0.59            | v        | 10.1 |     |
| Cadmium             |                             | < 0.0782            | u     | < 0.0306        | u        | NC   |     |
| Chromium            |                             | 12                  | v     | 16              | v        | 7.1  |     |
| Hexavalent Chromium |                             | < 0.64              | u     | 3.4             | v        | NC   |     |
| Cobalt              |                             | 4.9                 | v     | < 0.2828        | u        | NC   |     |
| Cyanide             |                             | 0.04                | J     | < 2.2626        | u        | NC   |     |
| Iron                |                             | 14000               | v     | 9100            | v        | 10.6 |     |
| Lead                |                             | 4.3                 | v     | 3.6             | v        | 4.4  |     |
| Manganese           |                             | 570                 | v     | 490             | v        | 3.8  |     |
| Mercury             |                             | 0.044               | v     | 0.035           | J        | 5.7  |     |

Table A-3  
Field Duplicate Summary  
SWMU 10 Investigation Report  
Western Refining Southwest, Inc. - Gallup Refinery

|  | Parameter | SWMU 10-22 (2-2.5') |   | DUP01           |   |      |
|--|-----------|---------------------|---|-----------------|---|------|
|  |           | 1605943-004         |   | 1605943-006     |   | RPD  |
|  |           | Sample Result       |   | Field Duplicate |   | %    |
|  | Nickel    | 8.6                 | v | 5.6             | v | 10.6 |
|  | Selenium  | < 1.346             | u | < 1.6905        | u | NC   |
|  | Silver    | < 0.0391            | u | < 0.0328        | u | NC   |
|  | Vanadium  | 18                  | v | 21              | v | 3.8  |
|  | Zinc      | 20                  | v | 24              | v | 4.5  |

**Notes:**

RPD = Relative percent difference;  $[(\text{difference})/(\text{average})] * 100$

NC = Not calculated; RPD values were not calculated for non-detects

ug/kg-dry = micrograms per kilogram dry

mg/kg-dry = milligrams per kilogram

bold value = Field Duplicate RPD Outlier

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                  | Parameter                      | SWMU 10-25 (2-2.5') |   | SWMU 10 DUP01   |   |      |
|------------------|--------------------------------|---------------------|---|-----------------|---|------|
|                  |                                | 1609B57-007         |   | 1609B57-010     |   | RPD  |
|                  |                                | Sample Result       |   | Field Duplicate |   | %    |
| TPH (mg/kg-dry)  | Gasoline Range Organics (GRO)  | 730                 | v | 510             | Z | 8.9  |
|                  | Diesel Range Organics (DRO)    | 790                 | v | 3400            | v | 31.1 |
|                  | Motor Oil Range Organics (MRO) | 100                 | v | < 560           | u | NC   |
| VOCs (mg/kg-dry) | 1,1,1,2-Tetrachloroethane      | < 0.2042            | u | < 0.18          | u | NC   |
|                  | 1,1,1-Trichloroethane          | < 0.1302            | u | < 0.11          | u | NC   |
|                  | 1,1,2,2-Tetrachloroethane      | < 0.3458            | u | < 0.3           | u | NC   |
|                  | 1,1,2-Trichloroethane          | < 0.2515            | u | < 0.22          | u | NC   |
|                  | 1,1-Dichloroethane             | < 0.1153            | u | < 0.1           | u | NC   |
|                  | 1,1-Dichloroethene             | < 0.6989            | u | < 0.61          | u | NC   |
|                  | 1,1-Dichloropropene            | < 0.1693            | u | < 0.15          | u | NC   |
|                  | 1,2,3-Trichlorobenzene         | < 0.3193            | u | < 0.28          | u | NC   |
|                  | 1,2,3-Trichloropropane         | < 0.3691            | u | < 0.32          | u | NC   |
|                  | 1,2,4-Trichlorobenzene         | < 0.2282            | u | < 0.2           | u | NC   |
|                  | 1,2,4-Trimethylbenzene         | 5.8                 | v | 5.8             | Z | 0.0  |
|                  | 1,2-Dibromo-3-chloropropane    | < 0.6536            | u | < 0.57          | u | NC   |
|                  | 1,2-Dibromoethane (EDB)        | < 0.1518            | u | < 0.13          | u | NC   |
|                  | 1,2-Dichlorobenzene            | < 0.1862            | u | < 0.16          | u | NC   |
|                  | 1,2-Dichloroethane (EDC)       | < 0.5564            | u | < 0.48          | u | NC   |
|                  | 1,2-Dichloropropane            | < 0.179             | u | < 0.16          | u | NC   |
|                  | 1,3,5-Trimethylbenzene         | 2.6                 | v | 2.5             | Z | 1.0  |
|                  | 1,3-Dichlorobenzene            | < 0.175             | u | < 0.15          | u | NC   |
|                  | 1,3-Dichloropropane            | < 0.2421            | u | < 0.21          | u | NC   |
|                  | 1,4-Dichlorobenzene            | < 0.2644            | u | < 0.23          | u | NC   |
|                  | 1-Methylnaphthalene            | 13                  | v | 13              | Z | 0.0  |
|                  | 2,2-Dichloropropane            | < 0.1222            | u | < 0.11          | u | NC   |
|                  | 2-Butanone                     | < 1.2194            | u | < 1.1           | u | NC   |
|                  | 2-Chlorotoluene                | < 0.1574            | u | < 0.14          | u | NC   |
|                  | 2-Hexanone                     | < 1.1611            | u | < 1             | u | NC   |
|                  | 2-Methylnaphthalene            | 16                  | v | 18              | Z | 2.9  |
|                  | 4-Chlorotoluene                | < 0.1887            | u | < 0.16          | u | NC   |
|                  | 4-Isopropyltoluene             | 0.61                | J | 0.59            | J | 0.8  |
|                  | 4-Methyl-2-pentanone           | < 0.6219            | u | < 0.54          | u | NC   |
|                  | Acetone                        | < 2.7608            | u | < 2.4           | u | NC   |
|                  | Benzene                        | < 0.8538            | u | < 0.74          | u | NC   |
|                  | Bromobenzene                   | < 0.172             | u | < 0.15          | u | NC   |
|                  | Bromodichloromethane           | < 0.1243            | u | < 0.11          | u | NC   |
|                  | Bromoform                      | < 0.2599            | u | < 0.23          | u | NC   |
|                  | Bromomethane                   | < 0.7861            | u | 0.69            | J | NC   |
|                  | Carbon disulfide               | < 0.7045            | u | < 0.61          | u | NC   |
|                  | Carbon tetrachloride           | < 0.1401            | u | < 0.12          | u | NC   |
|                  | Chlorobenzene                  | < 0.1737            | u | < 0.15          | u | NC   |
|                  | Chloroethane                   | < 0.4259            | u | < 0.37          | u | NC   |
|                  | Chloroform                     | < 0.161             | u | < 0.14          | u | NC   |
|                  | Chloromethane                  | < 0.1899            | u | < 0.17          | u | NC   |
|                  | cis-1,2-DCE                    | < 0.1241            | u | < 0.11          | u | NC   |
|                  | cis-1,3-Dichloropropene        | < 0.1967            | u | < 0.17          | u | NC   |

Table A-3  
Field Duplicate Summary  
SWMU 10 Investigation Report  
Western Refining Southwest, Inc. - Gallup Refinery

|                            | Parameter                      | SWMU 10-25 (2-2.5')    |          | SWMU 10 DUP01   |          |      |    |
|----------------------------|--------------------------------|------------------------|----------|-----------------|----------|------|----|
|                            |                                | 1609B57-007            |          | 1609B57-010     |          | RPD  |    |
|                            |                                | Sample Result          |          | Field Duplicate |          | %    |    |
|                            | Dibromochloromethane           | < 0.1928               | u        | < 0.17          | u        | NC   |    |
|                            | Dibromomethane                 | < 0.1848               | u        | < 0.16          | u        | NC   |    |
|                            | Dichlorodifluoromethane        | < 0.6602               | u        | < 0.57          | u        | NC   |    |
|                            | Ethylbenzene                   | 1                      | J        | 0.99            | J        | 0.3  |    |
|                            | Hexachlorobutadiene            | < 0.2609               | u        | < 0.23          | u        | NC   |    |
|                            | Isopropylbenzene               | 0.54                   | J        | 0.5             | J        | 1.9  |    |
|                            | Methyl tert-butyl ether (MTBE) | < 0.6706               | u        | < 0.58          | u        | NC   |    |
|                            | Methylene chloride             | < 0.6153               | u        | < 0.54          | u        | NC   |    |
|                            | Naphthalene                    | 3.8                    | J        | 4.1             | Z        | 1.9  |    |
|                            | n-Butylbenzene                 | 0.95                   | J        | 0.87            | J        | 2.2  |    |
|                            | n-Propylbenzene                | 0.93                   | J        | 0.98            | J        | 1.3  |    |
|                            | sec-Butylbenzene               | 0.74                   | J        | 0.67            | J        | 2.5  |    |
|                            | Styrene                        | < 0.1905               | u        | < 0.17          | u        | NC   |    |
|                            | tert-Butylbenzene              | < 0.1769               | u        | < 0.15          | u        | NC   |    |
|                            | Tetrachloroethene (PCE)        | < 0.177                | u        | < 0.15          | u        | NC   |    |
|                            | Toluene                        | 0.29                   | J        | 0.2             | J        | 9.2  |    |
|                            | trans-1,2-DCE                  | < 0.5973               | u        | < 0.52          | u        | NC   |    |
|                            | trans-1,3-Dichloropropene      | < 0.3123               | u        | < 0.27          | u        | NC   |    |
|                            | Trichloroethene (TCE)          | < 0.2288               | u        | < 0.2           | u        | NC   |    |
|                            | Trichlorofluoromethane         | < 0.1596               | u        | < 0.14          | u        | NC   |    |
|                            | Vinyl chloride                 | < 0.1745               | u        | < 0.15          | u        | NC   |    |
|                            | Xylenes, Total                 | 10                     | v        | 9.9             | Z        | 0.3  |    |
|                            | SVOCs (mg/kg-dry)              | 1,2,4-Trichlorobenzene | < 0.1232 | u               | < 1.1742 | u    | NC |
|                            |                                | 1,2-Dichlorobenzene    | < 0.0872 | u               | < 0.8314 | u    | NC |
| 1,3-Dichlorobenzene        |                                | < 0.088                | u        | < 0.839         | u        | NC   |    |
| 1,4-Dichlorobenzene        |                                | < 0.0963               | u        | < 0.9177        | u        | NC   |    |
| 1-Methylnaphthalene        |                                | 0.82                   | v        | 14              | Z        | 44.5 |    |
| 2,4,5-Trichlorophenol      |                                | < 0.114                | u        | < 1.0863        | u        | NC   |    |
| 2,4,6-Trichlorophenol      |                                | < 0.0945               | u        | < 0.9006        | u        | NC   |    |
| 2,4-Dichlorophenol         |                                | < 0.1063               | u        | < 1.0126        | u        | NC   |    |
| 2,4-Dimethylphenol         |                                | 0.17                   | J        | < 1.1788        | u        | NC   |    |
| 2,4-Dinitrophenol          |                                | < 0.0755               | u        | < 0.72          | u        | NC   |    |
| 2,4-Dinitrotoluene         |                                | < 0.1017               | u        | < 0.9694        | u        | NC   |    |
| 2,6-Dinitrotoluene         |                                | < 0.1205               | u        | < 1.1486        | u        | NC   |    |
| 2-Chloronaphthalene        |                                | < 0.0897               | u        | < 0.8546        | u        | NC   |    |
| 2-Chlorophenol             |                                | < 0.0898               | u        | < 0.8557        | u        | NC   |    |
| 2-Methylnaphthalene        |                                | 1.4                    | v        | 24              | Z        | 44.5 |    |
| 2-Methylphenol (cresol,o-) |                                | 0.53                   | v        | < 0.9073        | u        | NC   |    |
| 2-Nitroaniline             |                                | < 0.1228               | u        | < 1.1699        | u        | NC   |    |
| 2-Nitrophenol              |                                | < 0.113                | u        | < 1.0769        | u        | NC   |    |
| 3,3'-Dichlorobenzidine     |                                | < 0.0839               | u        | < 0.7992        | u        | NC   |    |
| 3+4-Methylphenol           |                                | < 0.0825               | u        | < 0.7858        | u        | NC   |    |
| 3-Nitroaniline             |                                | < 0.1004               | u        | < 0.9569        | u        | NC   |    |
| 4,6-Dinitro-2-methylphenol |                                | < 0.0689               | u        | < 0.6564        | u        | NC   |    |
| 4-Bromophenyl phenyl ether |                                | < 0.1089               | u        | < 1.0377        | u        | NC   |    |
| 4-Chloro-3-methylphenol    |                                | < 0.1359               | u        | < 1.2951        | u        | NC   |    |

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                    | Parameter                   | SWMU 10-25 (2-2.5') |          | SWMU 10 DUP01   |    |      |
|--------------------|-----------------------------|---------------------|----------|-----------------|----|------|
|                    |                             | 1609B57-007         |          | 1609B57-010     |    | RPD  |
|                    |                             | Sample Result       |          | Field Duplicate |    | %    |
|                    | 4-Chloroaniline             | < 0.1238            | u        | < 1.1801        | u  | NC   |
|                    | 4-Chlorophenyl phenyl ether | < 0.1301            | u        | < 1.2402        | u  | NC   |
|                    | 4-Nitroaniline              | < 0.0803            | u        | < 0.7654        | u  | NC   |
|                    | 4-Nitrophenol               | < 0.0868            | u        | < 0.8269        | u  | NC   |
|                    | Acenaphthene                | < 0.0977            | u        | < 0.9308        | u  | NC   |
|                    | Acenaphthylene              | < 0.0927            | u        | < 0.8833        | u  | NC   |
|                    | Aniline                     | < 0.1076            | u        | < 1.0256        | u  | NC   |
|                    | Anthracene                  | < 0.0756            | u        | < 0.7205        | u  | NC   |
|                    | Azobenzene                  | < 0.1388            | u        | < 1.3227        | u  | NC   |
|                    | Benz(a)anthracene           | < 0.0981            | u        | < 0.9344        | u  | NC   |
|                    | Benzo(a)pyrene              | < 0.0862            | u        | < 0.8219        | u  | NC   |
|                    | Benzo(b)fluoranthene        | < 0.1029            | u        | < 0.9803        | u  | NC   |
|                    | Benzo(g,h,i)perylene        | < 0.1004            | u        | < 0.9571        | u  | NC   |
|                    | Benzo(k)fluoranthene        | < 0.1003            | u        | < 0.9562        | u  | NC   |
|                    | Benzoic acid                | 0.21                | J        | 2.7             | J  | 42.8 |
|                    | Benzyl alcohol              | < 0.0892            | u        | < 0.8497        | u  | NC   |
|                    | Bis(2-chloroethoxy)methane  | < 0.1236            | u        | < 1.1778        | u  | NC   |
|                    | Bis(2-chloroethyl)ether     | < 0.0837            | u        | < 0.7976        | u  | NC   |
|                    | Bis(2-chloroisopropyl)ether | < 0.1017            | u        | < 0.9692        | u  | NC   |
|                    | Bis(2-ethylhexyl)phthalate  | 0.22                | J        | < 0.8844        | u  | NC   |
|                    | Butyl benzyl phthalate      | < 0.1008            | u        | < 0.961         | u  | NC   |
|                    | Carbazole                   | < 0.0769            | u        | < 0.733         | u  | NC   |
|                    | Chrysene                    | < 0.097             | u        | < 0.9243        | u  | NC   |
|                    | Dibenz(a,h)anthracene       | < 0.0921            | u        | < 0.8777        | u  | NC   |
|                    | Dibenzofuran                | < 0.1146            | u        | < 1.0917        | u  | NC   |
|                    | Diethyl phthalate           | 0.28                | v        | < 1.1007        | u  | NC   |
|                    | Dimethyl phthalate          | < 0.1114            | u        | < 1.0618        | u  | NC   |
|                    | Di-n-butyl phthalate        | 0.35                | J        | < 0.8117        | u  | NC   |
|                    | Di-n-octyl phthalate        | < 0.0972            | u        | < 0.9259        | u  | NC   |
|                    | Fluoranthene                | < 0.0656            | u        | < 0.6252        | u  | NC   |
|                    | Fluorene                    | 0.19                | J        | 2.8             | Z  | 43.6 |
|                    | Hexachlorobenzene           | < 0.0898            | u        | < 0.8559        | u  | NC   |
|                    | Hexachlorobutadiene         | < 0.1284            | u        | < 1.2232        | u  | NC   |
|                    | Hexachlorocyclopentadiene   | < 0.1302            | u        | < 1.241         | u  | NC   |
|                    | Hexachloroethane            | < 0.0979            | u        | < 0.9325        | u  | NC   |
|                    | Indeno(1,2,3-cd)pyrene      | < 0.0889            | u        | < 0.8476        | u  | NC   |
|                    | Naphthalene                 | 0.28                | v        | < 1.2           | u  | NC   |
|                    | Nitrobenzene                | < 0.1175            | u        | 4.7             | Z  | NC   |
|                    | N-Nitrosodi-n-propylamine   | < 0.1095            | u        | < 1.1201        | u  | NC   |
|                    | N-Nitrosodiphenylamine      | < 0.1112            | u        | < 1.0432        | u  | NC   |
|                    | Pentachlorophenol           | < 0.0732            | u        | < 1.06          | u  | NC   |
|                    | Phenanthrene                | 0.39                | v        | < 0.6976        | u  | NC   |
|                    | Phenol                      | < 0.0858            | u        | 5.1             | Z  | NC   |
| Pyrene             | < 0.0861                    | u                   | < 0.8175 | u               | NC |      |
| Pyridine           | < 0.0903                    | u                   | < 0.8203 | u               | NC |      |
| Metals (mg/kg-dry) | Antimony                    | < 1.1611            | u        | < 0.8604        | u  | NC   |

Table A-3  
Field Duplicate Summary  
SWMU 10 Investigation Report  
Western Refining Southwest, Inc. - Gallup Refinery

|  | Parameter           | SWMU 10-25 (2-2.5') |   | SWMU 10 DUP01   |   |      |
|--|---------------------|---------------------|---|-----------------|---|------|
|  |                     | 1609B57-007         |   | 1609B57-010     |   | RPD  |
|  |                     | Sample Result       |   | Field Duplicate |   | %    |
|  | Arsenic             | < 1.0254            | u | < 0.9931        | u | NC   |
|  | Barium              | 96                  | v | 130             | v | 7.5  |
|  | Beryllium           | 0.58                | v | 0.69            | v | 4.3  |
|  | Cadmium             | < 0.0732            | u | < 0.0709        | u | NC   |
|  | Chromium            | 7.2                 | v | 7.7             | v | 1.7  |
|  | Hexavalent Chromium | <0.1                | u | 6.1             | v | NC   |
|  | Cobalt              | 0.12                | J | 0.063           | J | 15.6 |
|  | Cyanide             | 0.3                 | J | 0.43            | J | 8.9  |
|  | Iron                | 12000               | v | 13000           | v | 2.0  |
|  | Lead                | 2.1                 | v | 3.4             | v | 11.8 |
|  | Manganese           | 500                 | v | 970             | v | 16.0 |
|  | Mercury             | 0.0031              | J | 0.0053          | J | 13.1 |
|  | Nickel              | 5.7                 | v | 6.7             | v | 4.0  |
|  | Selenium            | < 2.0993            | u | < 2.0332        | u | NC   |
|  | Silver              | < 0.0722            | u | < 0.07          | u | NC   |
|  | Vanadium            | 15                  | v | 17              | v | 3.1  |
|  | Zinc                | 11                  | v | 16              | v | 9.3  |

**Notes:**

RPD = Relative percent difference; [(difference)/(average)]\* 100

NC = Not calculated; RPD values were not calculated for non-detects

ug/kg-dry = micrograms per kilogram dry

mg/kg-dry = milligrams per kilogram

bold value = Field Duplicate RPD Outlier

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|             | Parameter                      | SWMU 10-15-GW | SWMU 10 DUP01GW | RPD (%) |
|-------------|--------------------------------|---------------|-----------------|---------|
|             |                                | Sample Result | Field Duplicate |         |
| TPH (mg/l): | Gasoline Range Organics (GRO)  | 0.78          | 0.73            | 1.7     |
|             | Diesel Range Organics (DRO)    | 1             | 1.1             | 2.4     |
|             | Motor Oil Range Organics (MRO) | 5 U           | 5 U             | NC      |
| VOCs (ug/l) | 1,1,1,2-Tetrachloroethane      | 0.22094 U     | 0.11047 U       | NC      |
|             | 1,1,1-Trichloroethane          | 0.15544 U     | 0.07772 U       | NC      |
|             | 1,1,2,2-Tetrachloroethane      | 0.35915 U     | 0.17957 U       | NC      |
|             | 1,1,2-Trichloroethane          | 0.15763 U     | 0.07881 U       | NC      |
|             | 1,1-Dichloroethane             | 0.8 U         | 0.4 U           | NC      |
|             | 1,1-Dichloroethene             | 0.19865 U     | 0.09932 U       | NC      |
|             | 1,1-Dichloropropene            | 0.23043 U     | 0.11521 U       | NC      |
|             | 1,2,3-Trichlorobenzene         | 0.53278 U     | 0.26639 U       | NC      |
|             | 1,2,3-Trichloropropane         | 0.31673 U     | 0.15836 U       | NC      |
|             | 1,2,4-Trichlorobenzene         | 0.56547 U     | 0.28273 U       | NC      |
|             | 1,2,4-Trimethylbenzene         | 0.31893 U     | 0.15946 U       | NC      |
|             | 1,2-Dibromo-3-chloropropane    | 0.27209 U     | 0.13604 U       | NC      |
|             | 1,2-Dibromoethane (EDB)        | 0.26750 U     | 0.13375 U       | NC      |
|             | 1,2-Dichlorobenzene            | 0.23661 U     | 0.11831 U       | NC      |
|             | 1,2-Dichloroethane (EDC)       | 0.35297 U     | 0.17648 U       | NC      |
|             | 1,2-Dichloropropane            | 0.30239 U     | 0.15119 U       | NC      |
|             | 1,3,5-Trimethylbenzene         | 0.24576 U     | 0.12288 U       | NC      |
|             | 1,3-Dichlorobenzene            | 0.18698 U     | 0.09349 U       | NC      |
|             | 1,3-Dichloropropane            | 0.34440 U     | 0.17220 U       | NC      |
|             | 1,4-Dichlorobenzene            | 0.33168 U     | 0.16584 U       | NC      |
|             | 1-Methylnaphthalene            | 1.07678 U     | 0.53839 U       | NC      |
|             | 2,2-Dichloropropane            | 0.30330 U     | 0.15165 U       | NC      |
|             | Benzene                        | 0.51 J        | 0.4 J           | 6.0     |
|             | 2-Butanone                     | 2.9 J         | 2.9 J           | 0.0     |
|             | 2-Chlorotoluene                | 0.15813 U     | 0.07906 U       | NC      |
|             | 2-Hexanone                     | 0.95322 U     | 0.47661 U       | NC      |
|             | 2-Methylnaphthalene            | 1.18893 U     | 0.59446 U       | NC      |
|             | 4-Chlorotoluene                | 0.29868 U     | 0.14934 U       | NC      |
|             | 4-Isopropyltoluene             | 0.37722 U     | 0.18861 U       | NC      |
|             | 4-Methyl-2-pentanone           | 0.51323 U     | 0.25661 U       | NC      |
|             | Acetone                        | 9.7 J         | 16              | 12.3    |
|             | Bromobenzene                   | 0.21555 U     | 0.10777 U       | NC      |
|             | Bromodichloromethane           | 0.17935 U     | 0.08968 U       | NC      |
|             | Bromoform                      | 0.32470 U     | 0.16235 U       | NC      |
|             | Bromomethane                   | 2.32245 U     | 1.16123 U       | NC      |
|             | Carbon disulfide               | 1.34646 U     | 0.67323 U       | NC      |
|             | Carbon Tetrachloride           | 0.15607 U     | 0.07803 U       | NC      |
|             | Chlorobenzene                  | 0.18680 U     | 0.09340 U       | NC      |
|             | Chloroethane                   | 0.20937 U     | 0.10468 U       | NC      |
|             | Chloroform                     | 0.42972 U     | 0.21486 U       | NC      |
|             | Chloromethane                  | 0.34737 U     | 0.17368 U       | NC      |
|             | cis-1,2-DCE                    | 0.16120 U     | 0.08060 U       | NC      |
|             | cis-1,3-Dichloropropene        | 0.26617 U     | 0.13308 U       | NC      |
|             | Dibromochloromethane           | 0.19361 U     | 0.09681 U       | NC      |
|             | Dibromomethane                 | 0.46872 U     | 0.23436 U       | NC      |
|             | Dichlorodifluoromethane        | 1.38481 U     | 0.69241 U       | NC      |
|             | Ethylbenzene                   | 0.20214 U     | 0.10107 U       | NC      |
|             | Hexachlorobutadiene            | 0.50280 U     | 0.25140 U       | NC      |
|             | Isopropylbenzene               | 0.30395 U     | 0.15197 U       | NC      |
|             | Methyl tert-butyl ether (MTBE) | 150           | 150             | 0.0     |
|             | Methylene Chloride             | 0.72072 U     | 0.36036 U       | NC      |
|             | Naphthalene                    | 0.43587 U     | 0.21793 U       | NC      |
|             | n-Butylbenzene                 | 0.49093 U     | 0.24546 U       | NC      |
|             | n-Propylbenzene                | 0.32634 U     | 0.16317 U       | NC      |
|             | sec-Butylbenzene               | 0.45720 U     | 0.22860 U       | NC      |
|             | Styrene                        | 0.21131 U     | 0.10566 U       | NC      |
|             | tert-Butylbenzene              | 0.25893 U     | 0.12947 U       | NC      |
|             | Tetrachloroethene (PCE)        | 0.32095 U     | 0.16047 U       | NC      |



**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                    | Parameter                   | SWMU 10-15-GW | SWMU 10 DUP01GW | RPD (%) |
|--------------------|-----------------------------|---------------|-----------------|---------|
|                    |                             | Sample Result | Field Duplicate |         |
|                    | Toluene                     | 0.21693 U     | 0.10847 U       | NC      |
|                    | trans-1,2-DCE               | 0.18810 U     | 0.09405 U       | NC      |
|                    | trans-1,3-Dichloropropene   | 0.22617 U     | 0.11309 U       | NC      |
|                    | Trichloroethene (TCE)       | 0.32669 U     | 0.16334 U       | NC      |
|                    | Trichlorofluoromethane      | 0.25403 U     | 0.12701 U       | NC      |
|                    | Vinyl chloride              | 0.25082 U     | 0.12541 U       | NC      |
|                    | Xylenes, Total              | 0.56497 U     | 0.28248 U       | NC      |
| SVOCs (mg/kg-dry): | 1,2,4-Trichlorobenzene      | 1.99336 U     | 1.99336 U       | NC      |
|                    | 1,2-Dichlorobenzene         | 1.89877 U     | 1.89877 U       | NC      |
|                    | 1,3-Dichlorobenzene         | 1.69038 U     | 1.69038 U       | NC      |
|                    | 1,4-Dichlorobenzene         | 1.23419 U     | 1.23419 U       | NC      |
|                    | 1-Methylnaphthalene         | 1.80147 U     | 1.80147 U       | NC      |
|                    | 2,4,5-Trichlorophenol       | 1.61671 U     | 1.61671 U       | NC      |
|                    | 2,4,6-Trichlorophenol       | 1.25834 U     | 1.25834 U       | NC      |
|                    | 2,4-Dichlorophenol          | 1.39496 U     | 1.39496 U       | NC      |
|                    | 2,4-Dimethylphenol          | 1.85381 U     | 1.85381 U       | NC      |
|                    | 2,4-Dinitrophenol           | 1.06974 U     | 1.06974 U       | NC      |
|                    | 2,4-Dinitrotoluene          | 1.43393 U     | 1.43393 U       | NC      |
|                    | 2,6-Dinitrotoluene          | 1.49035 U     | 1.49035 U       | NC      |
|                    | 2-Chloronaphthalene         | 1.71573 U     | 1.71573 U       | NC      |
|                    | 2-Chlorophenol              | 1.20402 U     | 1.20402 U       | NC      |
|                    | 2-Methylnaphthalene         | 2.24599 U     | 2.24599 U       | NC      |
|                    | 2-Methylphenol              | 1.24572 U     | 1.24572 U       | NC      |
|                    | 2-Nitroaniline              | 1.79467 U     | 1.79467 U       | NC      |
|                    | 2-Nitrophenol               | 1.23136 U     | 1.23136 U       | NC      |
|                    | 3,3'-Dichlorobenzidine      | 2.60844 U     | 2.60844 U       | NC      |
|                    | 3+4-Methylphenol            | 1.47532 U     | 1.47532 U       | NC      |
|                    | 3-Nitroaniline              | 1.47977 U     | 1.47977 U       | NC      |
|                    | 4,6-Dinitro-2-methylphenol  | 1.36786 U     | 1.36786 U       | NC      |
|                    | 4-Bromophenyl phenyl ether  | 1.40872 U     | 1.40872 U       | NC      |
|                    | 4-Chloro-3-methylphenol     | 1.35070 U     | 1.35070 U       | NC      |
|                    | 4-Chloroaniline             | 1.87394 U     | 1.87394 U       | NC      |
|                    | 4-Chlorophenyl phenyl ether | 2.01318 U     | 2.01318 U       | NC      |
|                    | 4-Nitroaniline              | 1.24171 U     | 1.24171 U       | NC      |
|                    | 4-Nitrophenol               | 1.40498 U     | 1.40498 U       | NC      |
|                    | Acenaphthene                | 1.88470 U     | 1.88470 U       | NC      |
|                    | Acenaphthylene              | 1.86589 U     | 1.86589 U       | NC      |
|                    | Aniline                     | 1.54599 U     | 1.54599 U       | NC      |
|                    | Anthracene                  | 1.60486 U     | 1.60486 U       | NC      |
|                    | Azobenzene                  | 2.00531 U     | 2.00531 U       | NC      |
|                    | Benz(a)anthracene           | 2.50627 U     | 2.50627 U       | NC      |
|                    | Benzo(a)pyrene              | 2.72398 U     | 2.72398 U       | NC      |
|                    | Benzo(b)fluoranthene        | 2.40284 U     | 2.40284 U       | NC      |
|                    | Benzo(g,h,i)perylene        | 3.12517 U     | 3.12517 U       | NC      |
|                    | Benzo(k)fluoranthene        | 2.51488 U     | 2.51488 U       | NC      |
|                    | Benzoic acid                | 1.02557 U     | 1.02557 U       | NC      |
|                    | Benzyl alcohol              | 1.17187 U     | 1.17187 U       | NC      |
|                    | Bis(2-chloroethoxy)methane  | 1.80486 U     | 1.80486 U       | NC      |
|                    | Bis(2-chloroethyl)ether     | 1.77363 U     | 1.77363 U       | NC      |
|                    | Bis(2-chloroisopropyl)ether | 2.05946 U     | 2.05946 U       | NC      |
|                    | Bis(2-ethylhexyl)phthalate  | 3.32558 U     | 3.32558 U       | NC      |
|                    | Butyl benzyl phthalate      | 2.37012 U     | 2.37012 U       | NC      |
|                    | Carbazole                   | 1.49023 U     | 1.49023 U       | NC      |
|                    | Chrysene                    | 2.15442 U     | 2.15442 U       | NC      |
|                    | Dibenz(a,h)anthracene       | 3.26832 U     | 3.26832 U       | NC      |
|                    | Dibenzofuran                | 1.90194 U     | 1.90194 U       | NC      |
|                    | Diethyl phthalate           | 1.67338 U     | 1.67338 U       | NC      |
|                    | Dimethyl phthalate          | 1.99487 U     | 1.99487 U       | NC      |
|                    | Di-n-butyl phthalate        | 2.10312 U     | 2.10312 U       | NC      |
|                    | Di-n-octyl phthalate        | 1.93926 U     | 1.93926 U       | NC      |
|                    | Fluoranthene                | 1.47557 U     | 1.47557 U       | NC      |

**Table A-3**  
**Field Duplicate Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                          | Parameter                 | SWMU 10-15-GW | SWMU 10 DUP01GW | RPD (%)     |
|--------------------------|---------------------------|---------------|-----------------|-------------|
|                          |                           | Sample Result | Field Duplicate |             |
|                          | Fluorene                  | 1.67191 U     | 1.67191 U       | NC          |
|                          | Hexachlorobenzene         | 2.02231 U     | 2.02231 U       | NC          |
|                          | Hexachlorobutadiene       | 1.94908 U     | 1.94908 U       | NC          |
|                          | Hexachlorocyclopentadiene | 1.47091 U     | 1.47091 U       | NC          |
|                          | Hexachloroethane          | 1.61223 U     | 1.61223 U       | NC          |
|                          | Indeno(1,2,3-cd)pyrene    | 2.45952 U     | 2.45952 U       | NC          |
|                          | Isophorone                | 1.95246 U     | 1.95246 U       | NC          |
|                          | Naphthalene               | 1.83350 U     | 1.83350 U       | NC          |
|                          | Nitrobenzene              | 1.49388 U     | 1.49388 U       | NC          |
|                          | N-Nitrosodimethylamine    | 1.41684 U     | 1.41684 U       | NC          |
|                          | N-Nitrosodi-n-propylamine | 2.03650 U     | 2.03650 U       | NC          |
|                          | N-Nitrosodiphenylamine    | 2.46700 U     | 2.46700 U       | NC          |
|                          | Pentachlorophenol         | 1.13741 U     | 1.13741 U       | NC          |
|                          | Phenanthrene              | 1.99666 U     | 1.99666 U       | NC          |
|                          | Phenol                    | 1.08528 U     | 1.08528 U       | NC          |
|                          | Pyrene                    | 2.31812 U     | 2.31812 U       | NC          |
|                          | Pyridine                  | 1.67388 U     | 1.67388 U       | NC          |
| Water Quality (mg/l)     | Chloride                  | 2300          | 2400            | 1.1         |
|                          | Fluoride                  | 0.55          | 0.55            | 0.0         |
|                          | Sulfate                   | 440           | 470             | 1.6         |
|                          |                           |               |                 |             |
| Dissolved Metals (mg/l): | Antimony (d)              | 0.001429741 U | 0.001 U         | NC          |
|                          | Arsenic (d)               | 0.011 J       | 0.011 J         | 0.0         |
|                          | Barium (d)                | 0.22          | 0.22            | 0.0         |
|                          | Beryllium (d)             | 0.000312692 U | 0.00033 J       | NC          |
|                          | Cadmium (d)               | 0.000966229 U | 0.001 U         | NC          |
|                          | Chromium (d)              | 0.0094        | 0.0094          | 0.0         |
|                          | Cobalt (d)                | 0.023         | 0.023           | 0.0         |
|                          | Iron (d)                  | 0.61          | 2.2             | <b>28.3</b> |
|                          | Lead (d)                  | 0.0026 J      | 0.0039 J        | 10.0        |
|                          | Manganese (d)             | 2.4           | 2.5             | 1.0         |
|                          | Nickel (d)                | 0.65          | 0.66            | 0.4         |
|                          | Selenium (d)              | 0.02          | 0.023           | 3.5         |
|                          | Silver (d)                | 0.00121688 U  | 0.001 U         | NC          |
|                          | Vanadium (d)              | 0.0057 J      | 0.0076 J        | 7.1         |
|                          | Zinc                      | 0.15          | 0.036           | <b>30.6</b> |
|                          |                           |               |                 |             |
| Total Metals (mg/l):     | Antimony                  | 0.00215761 U  | 0.002 U         | NC          |
|                          | Arsenic                   | 0.012         | 0.012           | 0.0         |
|                          | Barium                    | 0.37          | 0.31            | 4.4         |
|                          | Beryllium                 | 0.0015 J      | 0.0014 J        | 1.7         |
|                          | Cadmium                   | 0.000966229 U | 0.001 U         | NC          |
|                          | Chromium                  | 0.0062        | 0.0059 J        | 1.2         |
|                          | Chromium, Hexavalent      | 0.0005 U      | 0.00061         | NC          |
|                          | Cobalt                    | 0.025         | 0.025           | 0.0         |
|                          | Cyanide                   | 0.055         | 0.051           | 1.9         |
|                          | Iron                      | 9.4           | 7.4             | 6.0         |
|                          | Lead                      | 0.014         | 0.012           | 3.8         |
|                          | Manganese                 | 2.9           | 3               | 0.8         |
|                          | Mercury                   | 0.0000589 U   | 0.0000589 U     | NC          |
|                          | Nickel                    | 0.64          | 0.65            | 0.4         |
|                          | Selenium                  | 0.019         | 0.022           | 3.7         |
|                          | Silver                    | 0.002112456 U | 0.002 U         | NC          |
|                          | Vanadium                  | 0.019 J       | 0.017 J         | 2.8         |

**Notes:**

RPD = Relative percent difference; [(difference)/(average)]\* 100

NC = Not calculated; RPD values were not calculated for non-detects

ug/l = micrograms per liter

mg/l = milligrams per liter

bold value = Field Duplicate RPD Outlier

**Table A-4**  
**Completeness Summary**  
**SWMU 10 Investigation Report**  
**Western Refining Southwest, Inc. - Gallup Refinery**

|                                 | Parameter                      | Total Number of Results | Number of Usable Results | Percent Technical Compliance |
|---------------------------------|--------------------------------|-------------------------|--------------------------|------------------------------|
| <b>TPH :</b>                    | Diesel Range Organics (DRO)    | 100                     | 100                      | 100                          |
|                                 | Motor Oil Range Organics (MRO) | 100                     | 100                      | 100                          |
|                                 | Gasoline Range Organics (GRO)  | 100                     | 100                      | 100                          |
| <b>VOCs:</b>                    | All VOC Analytes               | 111                     | 111                      | 100                          |
| <b>SVOCs:</b>                   | All SVOC Analytes              | 100                     | 100                      | 100                          |
| <b>Metals (total):</b>          | Antimony                       | 98                      | 98                       | 100                          |
|                                 | Arsenic                        | 98                      | 98                       | 100                          |
|                                 | Barium                         | 98                      | 98                       | 100                          |
|                                 | Beryllium                      | 98                      | 98                       | 100                          |
|                                 | Cadmium                        | 98                      | 98                       | 100                          |
|                                 | Chromium                       | 98                      | 98                       | 100                          |
|                                 | Cobalt                         | 98                      | 98                       | 100                          |
|                                 | Cyanide                        | 98                      | 98                       | 100                          |
|                                 | Hexavalent Chromium            | 98                      | 98                       | 100                          |
|                                 | Iron                           | 98                      | 98                       | 100                          |
|                                 | Lead                           | 98                      | 98                       | 100                          |
|                                 | Manganese                      | 98                      | 98                       | 100                          |
|                                 | Mercury                        | 98                      | 98                       | 100                          |
|                                 | Nickel                         | 98                      | 98                       | 100                          |
|                                 | Selenium                       | 98                      | 98                       | 100                          |
|                                 | Silver                         | 98                      | 98                       | 100                          |
|                                 | Vanadium                       | 98                      | 98                       | 100                          |
|                                 | Zinc                           | 98                      | 98                       | 100                          |
| <b>Metals (mg/l dissolved):</b> | Antimony                       | 10                      | 10                       | 100                          |
|                                 | Arsenic                        | 10                      | 10                       | 100                          |
|                                 | Barium                         | 10                      | 10                       | 100                          |
|                                 | Beryllium                      | 10                      | 10                       | 100                          |
|                                 | Cadmium                        | 10                      | 10                       | 100                          |
|                                 | Chromium                       | 10                      | 10                       | 100                          |
|                                 | Cobalt                         | 10                      | 10                       | 100                          |
|                                 | Iron                           | 10                      | 10                       | 100                          |
|                                 | Lead                           | 10                      | 10                       | 100                          |
|                                 | Manganese                      | 10                      | 10                       | 100                          |
|                                 | Nickel                         | 10                      | 10                       | 100                          |
|                                 | Selenium                       | 10                      | 10                       | 100                          |
|                                 | Silver                         | 10                      | 10                       | 100                          |
|                                 | Vanadium                       | 10                      | 10                       | 100                          |
|                                 | Zinc                           | 10                      | 10                       | 100                          |
| <b>Water Quality (mg/l)</b>     | Chloride                       | 10                      | 10                       | 100                          |
|                                 | Fluoride                       | 10                      | 10                       | 100                          |
|                                 | Sulfate                        | 10                      | 10                       | 100                          |

**Notes:**

Number of samples used in completeness calculations includes soil samples, groundwater samples, soil and groundwater field duplicates, equipment rinsate, and field blanks.

Percent Technical Compliance = (Number of usable results / Number of reported results) \* 100

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## **Appendix F**

### **Historical Analyses**

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TABLE 2  
Ciniza Refinery Metals Results for Wastewater samples  
Taken During July and August, 1980.

| Metal    | New Well<br>Raw Water<br>7/23/80<br>0830 hrs. | Pond 3 Inlet<br>/Softener Waste<br>8/11/80<br>1000 hrs. | Cooling Water<br>Tower Blowdown<br>7/23/80<br>0830 hrs. | API Separator<br>Overflow<br>7/17/80<br>1330 hrs. | API Separator<br>Overflow<br>7/19/80<br>1300 hrs. | API Separator<br>Overflow<br>7/23/80<br>0830 hrs. |
|----------|---|---|---|---|---|---|
| Arsenic  | 0.003   | 0.031   | 0.013   | 0.004   | 0.005   | 0.015   |
| Barium   | 0.014   | 0.068   | 0.022   | 0.22  | 0.094   | 0.105   |
| Cadmium  | <0.001  | <0.001  | <0.001  | <0.001  | <0.001  | <0.001  |
| Chromium | <0.001  | 0.026   | 13.   | 0.91  | 0.64  | 1.2   |
| Lead     | <0.001  | <0.001  | 0.001   | <0.001  | <0.001  | <0.001  |
| Selenium | <0.001  | 0.097   | 0.025   | 0.015   | 0.018   | 0.024   |
| Silver   | <0.001  | 0.002   | 0.010   | 0.006   | 0.012   | 0.005   |

**CEC** Controls for Environmental Pollution, Inc.  
P.O. BOX 5351 • Santa Fe, New Mexico 87502

IN STATE 505/982-9841  
OUT OF STATE 800/545-2180

**REPORT OF ANALYSIS**

LAB # 84-03-404

**LE IDENTIFICATION**  
Seperator Sludge

**DATE COLLECTED**  
03/13/84

**TYPE OF ANALYSIS**

**mg/liter**

Silver

<0.01

Arsenic

<0.01

Barium

0.4

Cadmium

<0.001

Chromium

0.036

Chromium, Hexavalent

<0.01

Chromium

10

Mercury

<0.0004

Oil and Grease

26,000

Lead

<0.001

Lead

340

pH

9.30

Selenium

<0.01

Total Organic Carbon

>400,000

(ug/gram)

(ug/gram)

(ug/gram)

(units)

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**Appendix G**  
**Site-Specific Dilution/Attenuation Factor**  
**Calculations**

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### Calculation of Site-Specific Dilution/Attenuation Factor (DAF)

The DAF value was calculated using equation 33 from NMED's *Risk Assessment Guidance for Site Investigations and Remediation* (July, 2015).

$$DAF = 1 + \left( \frac{K * i * D}{I * L} \right) \quad DAF = 1 + \left( \frac{38.96 * 0.027 * 10.58}{0.000242 * 100} \right) = 295$$

Where:

$$D = (0.0112 * L^2)^{0.5} + D_a \left( 1 - \exp \left[ \frac{-L * I}{K * i * D_a} \right] \right)$$

$$D = (0.0112 * 100^2)^{0.5} + 12.5 \left( 1 - \exp \left[ \frac{-100 * 0.000242}{38.96 * 0.027 * 12.5} \right] \right) = 10.58 \text{ m}$$

K = Aquifer hydraulic conductivity (m/yr)  
i = Hydraulic gradient (m/m)  
D = Mixing zone depth (m)  
I = Infiltration rate (m/yr)  
L = Source length parallel to ground water flow (m)  
D<sub>a</sub> = Aquifer thickness (m)

Derivation of site-specific values:

K = 38.96 m/yr as determined from a slug test at well MW-1 (completed in Sonsela Sandstone aquifer) and originally presented in Appendix C of the Discharge Plan Application for Giant Refining Company, Ciniza Refinery, dated November 21, 1985.

i = 0.013 m/m as measured during Fall 2015 ground water sampling event (potentiometric surface map of Sonsela aquifer attached)

D = 10.58 m [lower of aquifer thickness (12.5 m measured at nearest well that penetrates Sonsela aquifer – OW-12) or calculated mixing zone depth (10.58m)]

I = 0.000182 m/yr derivation using EPA's HELP model as described below

L = 100 m – conservative source area length measured across Aeration Basin and surrounding soils with potential impact

D<sub>a</sub> = 12.5 m - saturated thickness measured at nearest well that penetrates Sonsela aquifer – OW-12 during August 2015 ground water sampling event

### Calculation of Infiltration Rate

Pursuant to EPA's *Soil Screening Guidance: User's Guide* (Second Edition, July 1996), infiltration rates can be calculated either of two ways: (1) assume that infiltration rate is



equivalent to recharge, or (2) use the EPA HELP model to estimate infiltration. Because the Gallup site is located in an area with low annual rainfall rates and high potential evapotranspiration rates, method 1 is not representative of site conditions. That is to say that it is unreasonable to assume that infiltration is equal to recharge.

EPA's HELP model was used to calculate the site-specific infiltration rate. Site-specific meteorological data was obtained from the Western Regional Climate Center and New Mexico State University, which operates a nearby weather station (Thoreau 5 ENE) as part of the NWS Cooperator Climate Stations. The weather station is located east northeast of Thoreau, New Mexico on HWY 371 and is approximately 16 miles south southeast of the Gallup Refinery.

Data obtained from the Thoreau 5 ENE station includes mean monthly temperature and average monthly precipitation. The average wind speed (11.3 km/hr) was obtained from the Western Regional Climate Center, as measured at the Gallup, NM airport. Daily solar radiation and quarterly relative humidity values were based on measurements from Albuquerque, NM. This data was obtained from the National Oceanic and Atmospheric Administration (NOAA) and is included in the HELP model's Weather Generator module. A review of the monthly average weather conditions (temperature and precipitation) at Gallup and Albuquerque as shown in the table below indicates very similar conditions such that use of quarterly relative humidity and solar radiation from Albuquerque should be sufficient to estimate conditions at the Western Refining Gallup Refinery. The quarterly relative humidity values used are 48%, 30%, 45%, and 50% for the first, second, third, and fourth quarters, respectively.

The HELP model soil profile was constructed with two layers. The Sonsela Aquifer is considered to be the uppermost aquifer in the area (reference the May 2000 RCRA Post-Closure Care Part B Permit Application). The upper layer (Layer 1.1) represents the unconsolidated deposits that overlie the Chinle Group (Petrified Forest Formation) bedrock. Based on the soil borings logs prepared from the recent investigation of SWMU 10, the predominant lithology in this upper layer is silty clay and clay, with lesser amounts of sandy clay. There are relatively thin (several feet thick) discontinuous layers of clayey sand and sand. The soil type chosen in the model was loam with an average thickness of 3 meters. Loam would normally be expected contain a higher percentage of silt and sand than observed in the most of the soil borings and is considered to be a conservative estimate of actual field conditions.

The land surface was assumed to have a poor stand of grass, which is more conservative than the other possible option of bare soil. The slope of the land surface was calculated to be 4.0 % based on surveyed elevations across the extent of SWMU 10. Based on the selected soil type (loam) the model default value for porosity is 0.463, field capacity is 0.232, wilting point is 0.116, and saturated hydraulic conductivity is  $3.7\text{E-}4$  cm/sec. These model default values are taken from the US Department of Agriculture. Soil permeability tests conducted on soil samples collected beneath the nearby Land Treatment Unit indicated a permeability of  $1.9\text{E-}5$  cm/sec, thus the model default hydraulic conductivity of  $3.7\text{E-}4$  cm/sec is a conservative estimate (see attached page 2 of Appendix C from RCRA Post-Closure Care Permit Application).

The second layer (Layer 1.2) in the model represents the upper portion of the Chinle Group (i.e., Petrified Forest Formation), which overlies the Sonsela Aquifer. This upper portion of the Chinle Group consists of mudstone/claystone. The default soil type from

the HELP model chosen to represent the upper Chinle Group was moderately compacted clay. The model default value for porosity is 0.456, field capacity is 0.419, and wilting point is 0.332. The default porosity value of 0.456 is not believed to be representative of the claystone bedrock, thus a literature search was conducted to find more appropriate values. An average porosity of 0.165 was identified for the Jurassic Morrison Formation (claystone) at Long Park, Colorado (USDI, 1963). A site-specific hydraulic conductivity was used to be more representative of actual site conditions than the program default value. Yield tests, including slug tests and pumping tests have been performed at the refinery to estimate the hydraulic conductivity of the upper portion of the Chinle Group (see attached documentation). A slug test performed on July 3, 1984 in well OW-4 indicated a hydraulic conductivity of  $4.0\text{E-}7$  cm/sec. A pump test was performed in well OW-24 on February 20, 1985 and it yielded a hydraulic conductivity of  $2.5\text{E-}7$  cm/sec. The higher of these two conductivity values ( $4.0\text{E-}7$  cm/sec) was utilized as a site-specific hydraulic conductivity for the lower layer (i.e., Chinle Group – mudstone/sandstone). The thickness of this layer was estimated at 24.4 meters from boring log for OW-12.

Using the model's synthetic weather generator and the aforementioned inputs, the model was run for a 40 year period to simulate potential infiltration (percolation or leakage through Layer 1.2). The model output is enclosed, showing the annual values. Over the modeled 40 year period, the average annual infiltration was 0.000182 meters/year. This average annual infiltration was used in the aforementioned calculation of the site-specific DAF value.

#### References

USDI, 1963, Porosity and Bulk Density of Sedimentary Rocks, United States Department of the Interior, Geological Survey Bulletin 1144-E, p.55.

# ***Project : SWMU 10***

*Location Specific Infiltration Rates*

***Model : HELP***

*An US EPA model for predicting landfill hydrologic processes and testing of effectiveness of landfill designs*

***Author : Your title Scott Crouch***

***Client : Title Ed Riege***

***Location : Albuquerque***

***11/11/2016***

# 1. Profile. EPA profile1

## Model Settings

[HELP] Case Settings

| Parameter                 | Value            | Units |
|---------------------------|------------------|-------|
| Runoff Method             | Model calculated | (-)   |
| Initial Moisture Settings | Model calculated | (-)   |

[HELP] Surface Water Settings

| Parameter        | Value               | Units |
|------------------|---------------------|-------|
| Runoff Area      | 100                 | (%%)  |
| Vegetation Class | Poor stand of grass | (-)   |

## Profile Structure

| Layer  | Top ( m) | Bottom ( m) | Thickness ( m) |
|--|----------|-------------|----------------|
|  Alluvium       | 0.0000   | -3.0000     | 3.0000         |
|  Chinle Bedrock | -2.9995  | -27.4000    | 24.4005        |

### 1.1. Layer. Alluvium

Top Slope Length: 100.0000  
Bottom Slope Length: 0.0000  
Top Slope: 4.0000  
Bottom Slope : 0.0000

[HELP] Vertical Perc. Layer Parameters

| Parameter             | Value  | Units     |
|-----------------------|--------|-----------|
| total porosity        | 0.463  | (vol/vol) |
| field capacity        | 0.232  | (vol/vol) |
| wilting point         | 0.116  | (vol/vol) |
| sat.hydr.conductivity | 3.7E-4 | (cm/sec)  |
| subsurface inflow     | 0      | (mm/year) |

### 1.2. Layer. Chinle Bedrock

Top Slope Length: 0.0000  
Bottom Slope Length: 0.0000  
Top Slope: 0.0000  
Bottom Slope : 0.0000

[HELP] Barrier Soil Liner Parameters

| Parameter             | Value  | Units     |
|-----------------------|--------|-----------|
| total porosity        | 0.165  | (vol/vol) |
| field capacity        | 0.419  | (vol/vol) |
| wilting point         | 0.332  | (vol/vol) |
| sat.hydr.conductivity | 4.0E-7 | (cm/sec)  |
| subsurface inflow     | 0      | (mm/year) |

Annual Totals rate (m)

|  |                   |            |                    |                 |
|--|-------------------|------------|--------------------|-----------------|
|  | Precipitation (m) | Runoff (m) | Evapotranspiration | Change in water |
|--|-------------------|------------|--------------------|-----------------|

|             |            |            | (m)        | storage (m) |
|-------------|------------|------------|------------|-------------|
| Year-1 (m)  | 1.9660E-01 | 0.0000E+00 | 1.9677E-01 | -1.7350E-04 |
| Year-2 (m)  | 3.0180E-01 | 9.1537E-04 | 3.0657E-01 | -5.7573E-03 |
| Year-3 (m)  | 2.3510E-01 | 0.0000E+00 | 2.4060E-01 | -5.7839E-03 |
| Year-4 (m)  | 2.3000E-01 | 2.5366E-03 | 2.3451E-01 | -7.1001E-03 |
| Year-5 (m)  | 2.5270E-01 | 0.0000E+00 | 2.4301E-01 | 9.6812E-03  |
| Year-6 (m)  | 1.5870E-01 | 5.4743E-06 | 1.4507E-01 | 1.3598E-02  |
| Year-7 (m)  | 1.8420E-01 | 0.0000E+00 | 1.8756E-01 | -3.4412E-03 |
| Year-8 (m)  | 2.5770E-01 | 1.2916E-05 | 2.4903E-01 | 8.6291E-03  |
| Year-9 (m)  | 1.9170E-01 | 6.6953E-04 | 2.1089E-01 | -1.9862E-02 |
| Year-10 (m) | 2.2820E-01 | 1.2865E-03 | 2.1390E-01 | 1.2470E-02  |
| Year-11 (m) | 2.3680E-01 | 1.5046E-03 | 2.5558E-01 | -2.5677E-02 |
| Year-12 (m) | 2.5940E-01 | 4.5220E-06 | 2.4044E-01 | 1.8948E-02  |
| Year-13 (m) | 1.8440E-01 | 1.1864E-05 | 1.9941E-01 | -1.5097E-02 |
| Year-14 (m) | 1.5860E-01 | 0.0000E+00 | 1.5141E-01 | 7.1785E-03  |
| Year-15 (m) | 2.4990E-01 | 0.0000E+00 | 2.5856E-01 | -8.6762E-03 |
| Year-16 (m) | 1.6700E-01 | 0.0000E+00 | 1.5472E-01 | 1.2275E-02  |
| Year-17 (m) | 1.3040E-01 | 0.0000E+00 | 1.1903E-01 | 1.1365E-02  |
| Year-18 (m) | 1.5020E-01 | 0.0000E+00 | 1.6500E-01 | -1.4840E-02 |
| Year-19 (m) | 2.0530E-01 | 0.0000E+00 | 1.9808E-01 | 7.1662E-03  |
| Year-20 (m) | 1.8180E-01 | 2.6583E-04 | 1.8840E-01 | -6.8733E-03 |
| Year-21 (m) | 2.3550E-01 | 6.9849E-06 | 2.2949E-01 | 5.9732E-03  |
| Year-22 (m) | 1.3750E-01 | 0.0000E+00 | 1.4770E-01 | -1.0197E-02 |
| Year-23 (m) | 2.3340E-01 | 3.9462E-03 | 2.2872E-01 | 7.1646E-04  |
| Year-24 (m) | 2.2170E-01 | 0.0000E+00 | 2.2689E-01 | -5.1976E-03 |
| Year-25 (m) | 1.4510E-01 | 5.4743E-06 | 1.3219E-01 | 1.2899E-02  |
| Year-26 (m) | 2.0130E-01 | 1.1479E-03 | 1.8065E-01 | 1.9493E-02  |
| Year-27 (m) | 2.3200E-01 | 0.0000E+00 | 2.2765E-01 | 4.3452E-03  |
| Year-28 (m) | 1.9260E-01 | 0.0000E+00 | 2.1651E-01 | -2.3992E-02 |
| Year-29 (m) | 2.3390E-01 | 1.2639E-06 | 2.3704E-01 | -3.1509E-03 |
| Year-30 (m) | 1.8890E-01 | 0.0000E+00 | 1.9277E-01 | -3.8754E-03 |
| Year-31 (m) | 2.4520E-01 | 1.2662E-04 | 2.3575E-01 | 9.3052E-03  |
| Year-32 (m) | 2.2790E-01 | 0.0000E+00 | 2.2550E-01 | 2.3827E-03  |
| Year-33 (m) | 3.1730E-01 | 5.2643E-03 | 3.1169E-01 | 2.3910E-04  |
| Year-34 (m) | 2.1170E-01 | 0.0000E+00 | 2.0234E-01 | 9.3490E-03  |
| Year-35 (m) | 2.7430E-01 | 9.5079E-04 | 2.7840E-01 | -5.0673E-03 |
| Year-36 (m) | 1.5090E-01 | 0.0000E+00 | 1.3599E-01 | 1.4901E-02  |
| Year-37 (m) | 2.1680E-01 | 1.5524E-04 | 2.5088E-01 | -3.4290E-02 |
| Year-38 (m) | 1.7490E-01 | 0.0000E+00 | 1.6844E-01 | 6.4508E-03  |
| Year-39 (m) | 2.1190E-01 | 1.0321E-03 | 1.8405E-01 | 2.6758E-02  |
| Year-40 (m) | 1.7540E-01 | 0.0000E+00 | 1.9942E-01 | -2.4131E-02 |
| Total (m)   | 8.3887E+00 | 1.9850E-02 | 8.3706E+00 | -9.0586E-03 |

(continued)

|             | Water budget balance (m) | Percolation or leakage through Layer 2 (m) | Average head on top of Layer 2 (m) |
|-------------|--------------------------|--|------------------------------------|
| Year-1 (m)  | -2.9526E-09              | 5.2603E-06                                 | 1.5356E-08                         |
| Year-2 (m)  | -4.5326E-09              | 6.6981E-05                                 | 1.0819E-07                         |
| Year-3 (m)  | -3.5308E-09              | 2.8523E-04                                 | 4.2733E-07                         |
| Year-4 (m)  | -3.4543E-09              | 5.7620E-05                                 | 1.4046E-07                         |
| Year-5 (m)  | -3.7952E-09              | 7.9856E-06                                 | 2.3292E-08                         |
| Year-6 (m)  | -2.3834E-09              | 2.3447E-05                                 | 2.0752E-08                         |
| Year-7 (m)  | -2.7664E-09              | 8.1705E-05                                 | 9.5470E-08                         |
| Year-8 (m)  | -3.8703E-09              | 3.1331E-05                                 | 9.1296E-08                         |
| Year-9 (m)  | -2.8790E-09              | 5.0396E-06                                 | 1.4646E-08                         |
| Year-10 (m) | -3.4272E-09              | 5.4365E-04                                 | 7.2411E-07                         |
| Year-11 (m) | -3.5564E-09              | 5.3959E-03                                 | 8.2432E-06                         |
| Year-12 (m) | -3.8958E-09              | 1.1126E-05                                 | 2.9653E-08                         |
| Year-13 (m) | -2.7694E-09              | 7.7901E-05                                 | 1.6158E-07                         |
| Year-14 (m) | -2.3819E-09              | 1.0173E-05                                 | 1.4647E-08                         |
| Year-15 (m) | -3.7531E-09              | 1.8462E-05                                 | 5.4615E-08                         |
| Year-16 (m) | -2.5081E-09              | 6.0052E-06                                 | 1.7436E-08                         |
| Year-17 (m) | -1.9584E-09              | 1.9551E-06                                 | 5.6773E-09                         |
| Year-18 (m) | -2.2558E-09              | 4.1669E-05                                 | 9.2553E-08                         |
| Year-19 (m) | -3.0833E-09              | 5.3926E-05                                 | 3.3493E-08                         |
| Year-20 (m) | -2.7304E-09              | 8.4890E-06                                 | 2.4744E-08                         |
| Year-21 (m) | -3.5369E-09              | 2.8949E-05                                 | 6.4945E-08                         |
| Year-22 (m) | -2.0650E-09              | 2.4329E-06                                 | 7.0344E-09                         |

|             |             |            |            |
|-------------|-------------|------------|------------|
| Year-23 (m) | -3.5053E-09 | 1.8935E-05 | 5.4598E-08 |
| Year-24 (m) | -3.3296E-09 | 6.2363E-06 | 1.8079E-08 |
| Year-25 (m) | -2.1792E-09 | 1.6165E-06 | 4.8035E-09 |
| Year-26 (m) | -3.0232E-09 | 5.5487E-06 | 1.6342E-08 |
| Year-27 (m) | -3.4843E-09 | 4.4125E-06 | 1.3844E-08 |
| Year-28 (m) | -2.8926E-09 | 8.3842E-05 | 1.3650E-07 |
| Year-29 (m) | -3.5128E-09 | 1.1656E-05 | 3.0979E-08 |
| Year-30 (m) | -2.8370E-09 | 7.6141E-06 | 2.2408E-08 |
| Year-31 (m) | -3.6825E-09 | 1.7414E-05 | 3.8111E-08 |
| Year-32 (m) | -3.4227E-09 | 1.5151E-05 | 4.4395E-08 |
| Year-33 (m) | -4.7654E-09 | 1.0183E-04 | 1.6100E-07 |
| Year-34 (m) | -3.1794E-09 | 6.1061E-06 | 1.8312E-08 |
| Year-35 (m) | -4.1196E-09 | 1.5303E-05 | 4.4619E-08 |
| Year-36 (m) | -2.2663E-09 | 4.3212E-06 | 8.5338E-09 |
| Year-37 (m) | -3.2560E-09 | 5.0865E-05 | 1.0231E-07 |
| Year-38 (m) | -2.6267E-09 | 1.0114E-05 | 2.9412E-08 |
| Year-39 (m) | -3.1824E-09 | 5.6373E-05 | 9.4730E-08 |
| Year-40 (m) | -2.6342E-09 | 1.0777E-04 | 2.2750E-07 |
| Total (m)   | -1.2599E-07 | 7.2903E-03 | 7.2903E-06 |

↓

$$\frac{7.2903E-03 \text{ meters}}{40 \text{ yrs}} = 1.82E-05 \text{ meters/yr}$$

# ALBUQUERQUE WSFO AIRPOR, NEW MEXICO (290234)

## Period of Record Monthly Climate Summary

Period of Record : 1/ 1/1914 to 12/31/2005

|                                   | Jan  | Feb  | Mar  | Apr  | May  | Jun  | Jul  | Aug  | Sep  | Oct  | Nov  | Dec  | Annual |
|-----------------------------------|------|------|------|------|------|------|------|------|------|------|------|------|--------|
| Average Max. Temperature (F)      | 47.2 | 53.2 | 60.6 | 70.0 | 79.4 | 89.3 | 91.7 | 88.9 | 82.4 | 71.0 | 56.9 | 47.7 | 69.9   |
| Average Min. Temperature (F)      | 23.4 | 27.8 | 33.0 | 40.8 | 50.1 | 59.2 | 64.6 | 62.9 | 56.0 | 44.0 | 31.4 | 24.4 | 43.1   |
| Average Total Precipitation (in.) | 0.37 | 0.40 | 0.52 | 0.54 | 0.63 | 0.61 | 1.38 | 1.46 | 0.96 | 0.88 | 0.46 | 0.46 | 8.67   |
| Average Total SnowFall (in.)      | 2.3  | 1.8  | 1.6  | 0.6  | 0.0  | 0.0  | 0.0  | 0.0  | 0.0  | 0.1  | 1.0  | 2.2  | 9.6    |
| Average Snow Depth (in.)          | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 0      |

Percent of possible observations for period of record.

Max. Temp.: 91.2% Min. Temp.: 91.2% Precipitation: 96.6% Snowfall: 80.1% Snow Depth: 80.1%

Check Station Metadata or Metadata graphics for more detail about data completeness.

# THOREAU 5 ENE, NEW MEXICO (298830)

## Period of Record Monthly Climate Summary

Period of Record : 1/ 1/1930 to 11/30/1992

|                                   | Jan  | Feb  | Mar  | Apr  | May  | Jun  | Jul  | Aug  | Sep  | Oct  | Nov  | Dec  | Annual |
|-----------------------------------|------|------|------|------|------|------|------|------|------|------|------|------|--------|
| Average Max. Temperature (F)      | 43.2 | 47.9 | 54.5 | 63.8 | 72.6 | 83.1 | 85.5 | 82.3 | 76.6 | 66.8 | 53.4 | 44.9 | 64.5   |
| Average Min. Temperature (F)      | 18.6 | 21.9 | 26.5 | 32.3 | 40.7 | 49.8 | 55.8 | 53.8 | 47.4 | 37.2 | 26.5 | 19.8 | 35.9   |
| Average Total Precipitation (in.) | 0.64 | 0.61 | 0.68 | 0.45 | 0.58 | 0.56 | 1.72 | 2.16 | 1.19 | 0.88 | 0.59 | 0.65 | 10.71  |
| Average Total SnowFall (in.)      | 8.3  | 5.6  | 5.3  | 1.1  | 0.4  | 0.0  | 0.0  | 0.0  | 0.0  | 1.4  | 3.1  | 8.1  | 33.3   |
| Average Snow Depth (in.)          | 1    | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 0    | 1    | 0      |

Percent of possible observations for period of record.

Max. Temp.: 46.1% Min. Temp.: 46.1% Precipitation: 68.4% Snowfall: 61.8% Snow Depth: 61.8%

Check Station Metadata or Metadata graphics for more detail about data completeness.

Western Regional Climate Center, [wrc@ddri.edu](mailto:wrc@ddri.edu)

# THOREAU 5 ENE, NEW MEXICO

## Station Metadata

From NCDC Station Historical Listing for NWS Cooperative Network  
 ObsTyp: t-Temperature-1, p-Daily precip-2, w-(blank), s-(blank), e-Evap-5  
 h-Hourly precip - 6 0.01" Universal, or - 7 0.10" Fisher-Porter  
 U - Observed, but beginning date is uncertain

| Count | Number   | Station Name        | Lat  | Long  | Elev  | Start | ObsTyp      | End     |
|-------|----------|---------------------|------|-------|-------|-------|-------------|---------|
|       | (Coop)   | (From NCDC listing) | ddmm | dddmm | ftx10 | yy mm | t p w s e h | yy mm   |
| 1050  | 298830-4 | THOREAU 6 ENE       | 3526 | 10808 | 0713  | 53 09 | U U         | 63 01   |
| 1051  | 298830-4 | THOREAU 6 ENE       | 3526 | 10809 | 0711  | 63 01 | U U         | 88 04   |
| 1052  | 298830-4 | THOREAU 5 ENE       | 3526 | 10809 | 0710  | 88 04 | 1 2         | 7 99 99 |

## Statistics by element

(From WRCC data archives)  
 Last Compiled on Apr.23, 2009.

Dates are format of YYYYMMDD. Numbers are total Number of observations

| STATION | START    | END      | PRECIP | SNWFL | SNWDP | TMAX  | TMIN  | TOBS  | EVAP | WNDM |
|---------|----------|----------|--------|-------|-------|-------|-------|-------|------|------|
| 298830  | 19300101 | 19921130 | 15723  | 14204 | 14192 | 10584 | 10584 | 10580 | 0    | 0    |

STATION - NCDC COOP Station number  
 START - First Date in record  
 END - Last Date in record (when last compiled)  
 PRECIP - Precipitation  
 SNWFL - Snowfall  
 SNWDP - Snow depth  
 TMAX - Daily Max. Temperature  
 TMIN - Daily Min. Temperature  
 TOBS - Temperature at Observation time  
 EVAP - Evaporation  
 WNDMV - Wind Movement

## Statistics by observation

(From WRCC data archives)  
 Last Compiled on Apr.23, 2009.

Dates are format of YYYYMMDD. Numbers represent one day and one day is considered present if



## Average Wind Speeds - MPH

Average wind speeds are based on the hourly data from 1996-2006 from automated stations at reporting airports (ASOS) unless otherwise noted.

For more information click here on [Western States](#), [Alaska](#) or [Hawaii and Pacific Islands](#)

[Arizona](#) [California](#) [Colorado](#) [Hawaii](#) [Idaho](#) [Montana](#) [Nevada](#) [New Mexico](#) [Oregon](#) [Utah](#) [Washington](#) [Wyoming](#)

### MONTANA

#### AVERAGE WIND SPEED - MPH

| STATION                 | ID             | Years | Jan  | Feb  | Mar  | Apr  | May  | Jun  | Jul  | Aug  | Sep  | Oct  | Nov  | Dec  | Ann  |
|-------------------------|----------------|-------|------|------|------|------|------|------|------|------|------|------|------|------|------|
| BAKER AIRPORT ASOS      | KBHK 1998-2006 |       | 10.4 | 10.5 | 12.2 | 12.5 | 12.7 | 11.7 | 10.7 | 10.8 | 10.0 | 10.3 | 10.5 | 10.7 | 11.1 |
| BILLINGS-LOGAN AP ASOS  | KBIL 1996-2006 |       | 12.8 | 12.2 | 11.0 | 10.5 | 10.4 | 10.0 | 9.3  | 9.2  | 9.6  | 10.3 | 12.0 | 13.0 | 10.9 |
| BOZEMAN AIRPORT ASOS    | KBZN 1996-2006 |       | 4.6  | 5.3  | 6.4  | 7.2  | 7.0  | 6.3  | 6.3  | 6.4  | 6.1  | 5.7  | 4.9  | 4.8  | 5.9  |
| BUTTE AIRPORT ASOS      | KBTM 1996-2006 |       | 4.7  | 5.0  | 6.8  | 7.3  | 7.7  | 7.3  | 6.9  | 6.6  | 6.2  | 6.2  | 5.2  | 4.9  | 6.2  |
| CUT BANK AIRPORT ASOS   | KCTB 1996-2006 |       | 14.6 | 12.8 | 13.3 | 12.8 | 13.4 | 12.3 | 11.2 | 10.2 | 11.4 | 13.0 | 13.6 | 14.6 | 12.8 |
| DILLON AIRPORT ASOS     | KDLN 1997-2006 |       | 10.2 | 9.7  | 10.3 | 10.0 | 9.6  | 8.4  | 8.0  | 8.0  | 8.7  | 9.0  | 9.3  | 9.2  | 9.2  |
| GLASGOW AIRPORT ASOS    | KGGW 1996-2006 |       | 9.3  | 9.3  | 11.7 | 12.4 | 12.6 | 11.4 | 10.7 | 11.0 | 10.5 | 10.3 | 9.3  | 9.4  | 10.6 |
| GLENDIVE AIRPORT AWOS   | KGDV 1996-2006 |       | 9.5  | 9.7  | 10.5 | 11.3 | 11.6 | 10.4 | 9.4  | 9.6  | 9.6  | 10.1 | 9.7  | 10.2 | 10.1 |
| GREAT FALLS AP ASOS     | KGTf 1996-2006 |       | 13.3 | 12.3 | 11.8 | 11.2 | 11.3 | 10.2 | 9.6  | 9.2  | 10.4 | 11.9 | 13.2 | 13.8 | 11.5 |
| GREAT FALLS-MALSTROM AF | KGFA 1996-2006 |       | 12.3 | 9.9  | 11.9 | 10.7 | 10.3 | 9.7  | 9.1  | 8.8  | 9.9  | 10.4 | 13.2 | 13.2 | 10.9 |
| HAVRE AIRPORT ASOS      | KHVR 1996-2006 |       | 9.9  | 9.6  | 10.7 | 11.0 | 11.6 | 10.6 | 9.8  | 9.5  | 9.8  | 9.8  | 10.6 | 11.0 | 10.3 |
| HELENA AIRPORT ASOS     | KHLN 1996-2006 |       | 5.8  | 6.3  | 7.8  | 8.3  | 8.4  | 8.2  | 7.4  | 6.6  | 6.7  | 6.6  | 5.8  | 6.0  | 7.0  |
| JORDAN AIRPORT ASOS     | KJDN 1996-2006 |       | 7.3  | 7.9  | 9.3  | 10.0 | 10.5 | 9.7  | 8.4  | 8.4  | 8.2  | 8.2  | 7.9  | 8.1  | 8.6  |
| KALISPELL AIRPORT ASOS  | KGPI 1996-2006 |       | 4.1  | 3.8  | 6.0  | 6.7  | 6.6  | 5.7  | 5.2  | 5.0  | 4.6  | 4.2  | 4.0  | 3.2  | 4.9  |
| LEWISTOWN AIRPORT ASOS  | KLWT 1996-2006 |       | 10.6 | 9.5  | 10.2 | 9.8  | 9.9  | 9.0  | 7.9  | 8.0  | 8.5  | 9.2  | 10.2 | 10.7 | 9.5  |
| LIVINGSTON AIRPORT ASOS | KLVM 1996-2006 |       | 19.8 | 17.4 | 16.2 | 14.0 | 13.1 | 11.9 | 11.0 | 11.2 | 12.7 | 14.6 | 18.4 | 20.4 | 15.2 |
| MILES CITY AP ASOS      | KMLS 1996-2006 |       | 8.8  | 9.4  | 10.6 | 11.2 | 11.3 | 10.5 | 9.9  | 9.7  | 9.7  | 9.7  | 9.3  | 9.3  | 9.9  |
| MISSOULA AIRPORT ASOS   | KMSO 1996-2006 |       | 3.2  | 3.7  | 5.4  | 6.1  | 6.0  | 6.1  | 5.7  | 5.2  | 4.3  | 3.9  | 3.5  | 3.7  | 4.7  |
| SIDNEY AIRPORT AWOS     | KSDY 1996-2006 |       | 8.9  | 9.0  | 9.5  | 10.2 | 10.4 | 9.0  | 7.7  | 7.9  | 8.2  | 8.8  | 8.7  | 9.4  | 9.0  |
| WOLF POINT AIRPORT ASOS | KOLF 1998-2006 |       | 7.1  | 7.6  | 9.4  | 10.3 | 10.4 | 8.9  | 8.2  | 8.3  | 7.6  | 7.7  | 7.3  | 7.6  | 8.3  |

### NEW MEXICO

#### AVERAGE WIND SPEED - MPH

| STATION                 | ID             | Years | Jan  | Feb  | Mar  | Apr  | May  | Jun  | Jul  | Aug  | Sep  | Oct  | Nov  | Dec  | Ann  |
|-------------------------|----------------|-------|------|------|------|------|------|------|------|------|------|------|------|------|------|
| ALAMOGORDO AIRPORT ASOS | KALM 1996-2006 |       | 5.1  | 6.3  | 7.1  | 7.9  | 7.1  | 6.9  | 6.1  | 5.3  | 5.2  | 5.2  | 5.0  | 5.0  | 6.0  |
| ALAMOGORDO-HOLLOMAN AFB | KHMN 1996-2006 |       | 8.5  | 9.7  | 10.6 | 11.8 | 10.8 | 10.6 | 9.8  | 9.1  | 8.8  | 8.5  | 8.1  | 8.3  | 9.6  |
| ALBUQUERQUE AP ASOS     | KABQ 1996-2006 |       | 7.0  | 8.2  | 9.3  | 11.1 | 10.0 | 10.0 | 8.7  | 8.3  | 8.0  | 7.9  | 7.2  | 6.9  | 8.5  |
| ALBUQUERQUE-DBLE EAGLE  | KAEG 1999-2006 |       | 7.1  | 7.9  | 9.0  | 10.6 | 9.5  | 8.6  | 7.0  | 6.2  | 7.0  | 6.5  | 6.5  | 6.1  | 7.7  |
| ARTESIA AIRPORT ASOS    | KATS 1997-2006 |       | 7.8  | 9.1  | 10.1 | 10.9 | 10.2 | 9.9  | 7.8  | 6.9  | 7.6  | 7.8  | 7.6  | 7.4  | 8.5  |
| CARLSBAD AIRPORT ASOS   | KCNM 1996-2006 |       | 9.2  | 9.8  | 10.9 | 11.4 | 10.4 | 9.9  | 8.5  | 7.7  | 8.2  | 8.5  | 8.4  | 8.8  | 9.3  |
| CLAYTON MUNI AP ASOS    | KCAO 1996-2006 |       | 11.9 | 12.7 | 13.4 | 14.6 | 13.4 | 13.0 | 11.7 | 10.8 | 11.8 | 12.1 | 12.1 | 12.0 | 12.4 |
| CLINES CORNERS          | KCQC 1998-2006 |       | 16.2 | 16.1 | 15.7 | 16.9 | 14.6 | 13.5 | 10.6 | 10.1 | 11.8 | 13.3 | 15.0 | 16.0 | 14.1 |
| CLOVIS AIRPORT AWOS     | KCVN 1996-2006 |       | 12.3 | 12.3 | 13.4 | 13.8 | 12.4 | 11.9 | 9.7  | 8.9  | 9.7  | 10.9 | 11.6 | 12.2 | 11.6 |
| CLOVIS-CANNON AFB       | KCVS 1996-2006 |       | 12.5 | 12.6 | 13.6 | 13.8 | 12.2 | 12.5 | 10.7 | 10.0 | 10.2 | 11.3 | 11.7 | 12.4 | 12.0 |
| DEMING AIRPORT ASOS     | KDMN 1996-2006 |       | 8.7  | 9.7  | 10.9 | 12.0 | 10.6 | 10.1 | 8.9  | 8.1  | 8.4  | 8.2  | 8.5  | 8.1  | 9.3  |
| FARMINGTON AIRPORT ASOS | KFMN 1996-2006 |       | 7.3  | 8.3  | 9.0  | 9.8  | 9.4  | 9.4  | 8.7  | 8.2  | 8.0  | 7.8  | 7.6  | 7.3  | 8.4  |
| GALLUP AIRPORT ASOS     | KGUP 1996-2006 |       | 5.7  | 6.9  | 7.8  | 10.0 | 9.0  | 8.8  | 6.9  | 6.0  | 6.5  | 6.1  | 5.6  | 5.3  | 7.0  |
| GRANTS-MILAN AP ASOS    | KGNT 1997-2006 |       | 7.8  | 8.8  | 9.6  | 10.9 | 10.0 | 9.8  | 8.1  | 7.2  | 7.9  | 8.4  | 8.0  | 7.6  | 8.7  |
| HOBBS AIRPORT AWOS      | KHOB 1996-2006 |       | 11.3 | 11.9 | 12.6 | 13.4 | 12.5 | 12.3 | 11.0 | 10.0 | 10.2 | 10.6 | 10.7 | 11.1 | 11.4 |
| LAS CRUCES AIRPORT AWOS | KLRU 2000-2006 |       | 6.4  | 7.5  | 8.8  | 10.1 | 8.7  | 8.2  | 6.8  | 6.0  | 6.2  | 6.1  | 6.4  | 6.0  | 7.3  |
| LAS VEGAS AIRPORT ASOS  | KLVS 1996-2006 |       | 10.9 | 12.2 | 12.5 | 14.3 | 12.4 | 11.8 | 10.0 | 9.2  | 10.9 | 10.8 | 11.0 | 10.9 | 11.4 |
| LOS ALAMOS AP AWOS      | KLAM 2005-2006 |       | 3.9  | 5.7  | 7.5  | 8.1  | 7.1  | 7.3  | 5.3  | 4.8  | 5.7  | 5.1  | 4.4  | 3.2  | 5.4  |
| RATON AIRPORT ASOS      | KRTN 1998-2006 |       | 8.9  | 9.4  | 10.4 | 12.2 | 10.8 | 10.2 | 8.4  | 8.1  | 8.6  | 9.0  | 8.6  | 8.5  | 9.4  |
| ROSWELL AIRPORT ASOS    | KROW 1996-2006 |       | 7.4  | 8.9  | 9.9  | 11.1 | 10.3 | 10.2 | 8.8  | 7.9  | 8.3  | 8.0  | 7.5  | 7.3  | 8.8  |
| RUIDOSO AIRPORT AWOS    | KSRR 1996-2006 |       | 8.8  | 9.6  | 10.0 | 11.6 | 10.0 | 8.4  | 5.9  | 5.3  | 6.4  | 7.4  | 7.9  | 8.7  | 8.3  |
| SANTA FE AIRPORT ASOS   | KSAF 1996-2006 |       | 8.9  | 9.5  | 9.9  | 11.2 | 10.6 | 10.5 | 9.2  | 8.8  | 8.8  | 9.1  | 8.7  | 8.5  | 9.5  |
| SILVER CITY AP AWOS     | KSVC 1999-2006 |       | 8.1  | 8.7  | 9.9  | 10.8 | 10.2 | 9.9  | 8.5  | 7.2  | 6.9  | 7.6  | 7.9  | 7.7  | 8.5  |
| TAOS AIRPORT AWOS       | KSKX 1996-2006 |       | 5.8  | 6.5  | 7.7  | 9.1  | 8.6  | 8.5  | 7.1  | 6.6  | 6.7  | 6.6  | 6.0  | 5.7  | 7.0  |
| TRUTH OR CONSEQ AP ASOS | KTCS 1996-2006 |       | 7.4  | 8.7  | 9.9  | 11.1 | 10.4 | 9.8  | 8.1  | 7.4  | 7.7  | 8.0  | 7.7  | 7.3  | 8.6  |
| TUCUMCARI AIRPORT ASOS  | KTCC 1999-2006 |       | 10.0 | 11.2 | 11.9 | 13.6 | 11.9 | 11.6 | 9.9  | 9.3  | 10.0 | 10.0 | 10.4 | 10.2 | 10.8 |

$$7.0 \text{ miles/hr} \times \frac{1 \text{ kilo.}}{0.62 \text{ miles}} = 11.3 \text{ K/hr}$$

SONSELA POTENTIOMETRIC SURFACE MAP  
WESTERN REFINING COMPANY  
GALLUP REFINERY

Fall 2015



**psi**  
4601 Ripley  
El Paso, Texas  
79922  
915-394-1317

1"=500'  
Project #: 06251104

**Figure 9**  
**Sonsela Water Elevation Map 2015**  
**WESTERN REFINING - GALLUP REFINERY**

Western Refining - Gallup Refinery  
92 Giant Crossing Road  
Gallup, New Mexico 87301  
Date: August 5, 2016 - Revision 1

APPENDIX C – DISCHARGE PLAN APPLICATION  
GIANT REFINING COMPANY  
CINIZA REFINERY

November 21, 1985

APPENDIX C  
AQUIFER-TEST DATA AND ANALYSES

IDENT.  
TEST PUMPING OF  
CHINLE SHALE

METHODOLOGY AND DESCRIPTION OF THE TEST

The test consisted of a 5 hour pumping period and a 2 hour recovery period. An air-driven piston pump capable of sustaining pumping rates as low as 10 gallons/hour (0.167 gpm) was used for the test. Water level measurements were taken with an electronic sounder. The well (OW-24) is located approximately 250 feet northwest of the land treatment facility and is completed within the Chinle shale. The lithologic and completion log of the well is attached (Figure F-2).

Pumping began at 1515 hours on February 20, 1985 at a rate of 10 gallons/hour. The produced water was very turbid. Clogging of the pump and pump lines necessitated continuous monitoring and adjustment of the discharge.

After 4 hours of pumping at 10 gallons/hour, the drawdown of the well appeared to stabilize at about 7 feet. The discharge rate was increased to 20 gallons/hour in order to more effectively stress the aquitard. After one hour of additional pumping a total drawdown of 12 feet was observed. However, this higher pumping rate increased the turbidity of the discharge and caused instability of the pumping rate. The lack of control of the discharge rate and the potential of damage to the pump forced the termination of the test after a total of 5 hours of pumping.

Water level recovery was observed for 100 minutes. At this time the water level had recovered to within 90% of the pre-pumping level.

TABLE F-1

Pump Test Data, OW-24

### PUMP TEST ANALYSIS

Field measurements are summarized in Table F-1. Due to the short pumping time and potential well-bore and gravel-pack effects, the final analysis was based on methods developed by Shafer, for low-conductivity materials.

Partial penetration effects were neglected in the analysis because the low pumping rates and the expected anisotropy of the aquitard would prevent significant vertical flow to the well bore. The low pumping rate was also designed to completely drain the gravel pack in the well to insure accurate recovery data.

A copy of Shafer's methodology is attached, and the data for his analysis is given in Table F-2. Figure F-1 is a plot of the recovery data, according to Shafer's methods. This Figure includes calculation of T and K for the Chinle shales.



$$TD = 58' - 4" = 58.33'$$

$$DW = 31.96'$$

# PUMPING DATA

Page 1 of 2

Pump C 55 26.37' of water

≈ 18 gal

## DATA SHEET FOR RECORDING PUMP TEST DATA

County: McKinley Co.

Observation well no.

Location: Cimiza Refinery 2/2/85

Pumped well no. OW-24

Average Q = \_\_\_\_\_ gpm r = \_\_\_\_\_ ft. r<sup>2</sup> = \_\_\_\_\_

| Date   | Hour | t (min) | t' (min) | t/t' | Depth to water | s (unadjusted) | Adjustment Δs | s' (adjusted) | gph (approx) | Remarks              |
|--------|------|---------|----------|------|----------------|----------------|---------------|---------------|--------------|----------------------|
| 2-20   | 1515 | 0       |          |      | 31'-11.5"      | 0              | Feet          |               | 40 gph       | STATIC LEVEL w/ PUMP |
|        | 1516 | 1       |          |      | 33-0           | 12.5           | 1.04          |               | 10 gph       | LARGE INITIAL SURGE  |
|        | 1517 | 2       |          |      | 33-1           | 13.5           | 1.13          |               |              | FLIGHT HOSE FILLING  |
|        | 1518 | 3       |          |      | 33-4           | 16.5           | 1.38          |               |              | Setting pump rate    |
|        | 1519 | 4       |          |      | 33-8 3/4       | 20.75          | 1.73          |               |              |                      |
|        | 1520 | 5       |          |      | 33-11 1/2      | 24.0           | 2.0           |               |              |                      |
|        | 1521 | 6       |          |      | 34-2           | 26.5           | 2.22          |               |              |                      |
|        | 1522 | 7       |          |      | 34-2 1/2       | 27.0           | 2.25          |               |              |                      |
| 1 min  | 1523 | 8       |          |      | 34-3           | 27.5           | 2.27          |               |              |                      |
|        | 1524 | 9       |          |      | 34-4           | 28.5           | 2.37          |               |              |                      |
|        | 1525 | 10      |          |      | 34-5           | 29.5           | 2.46          |               |              |                      |
|        | 1527 | 12      |          |      | 34-7 1/4       | 30.75          | 2.56          |               |              |                      |
|        | 1529 | 14      |          |      | 34-9           | 33.5           | 2.79          |               |              |                      |
| 2 min  | 1531 | 16      |          |      | 35-0           | 36.5           | 3.04          |               |              |                      |
|        | 1533 | 18      |          |      | 35-2 1/4       | 38.75          | 3.23          |               |              |                      |
|        | 1535 | 20      |          |      | 35-4 1/2       | 41.0           | 3.42          |               |              |                      |
|        | 1540 | 25      |          |      | 35-10 1/4      | 47.0           | 3.92          |               |              |                      |
|        | 1545 | 30      |          |      | 36-3           | 51.5           | 4.28          |               |              |                      |
|        | 1550 | 35      |          |      | 36-2 1/2       | 51.0           | 4.25          |               |              |                      |
| 5 min  | 1555 | 40      |          |      | 36-8           | 56.5           | 4.71          |               | 10 gph       | CLEARING PUMP Valve  |
|        | 1600 | 45      |          |      | 37-2 1/4       | 61.75          | 5.23          |               |              |                      |
|        | 1605 | 50      |          |      | 36-10 1/2      | 61.0           | 5.08          |               |              | clear pump valve     |
|        | 1610 | 55      |          |      | NR             | 70.75          |               |               |              |                      |
|        | 1615 | 60      |          |      | 37-10 1/4      | 82.75          | 5.90          |               |              | Pumped 10 gal        |
| 10 min | 1625 | 70      |          |      | 38-1 1/2       | 85.0           | 6.08          |               |              |                      |

# PUMPING DATA

Page 2 of 2

## DATA SHEET FOR RECORDING PUMP TEST DATA

County: \_\_\_\_\_

Observation well no. MW-24

Location: Ciniga Refinery

Pumped well no. MW-24

Average 0.10 gpm  $q =$  \_\_\_\_\_  $sc. r^2 =$  \_\_\_\_\_

| Date    | Hour | t (min) | t' (min) | t/t' | Depth to water | s // (unad-justed) | Adjust-ment Δs | s' (ad-justed) | gpm | Remarks  |
|---------|------|---------|----------|------|----------------|--------------------|----------------|----------------|-----|--|
| 2-20    | 1635 | 80      |          |      | 39-3 1/2       | 78                 | 6.33           |                | 10  |  |
|         | 1645 | 90      |          |      | 38-7           | 79.5               | 6.79           |                |     | Lift pump to 50'   |
|         | 1655 | 100     |          |      | 37-8 1/2       | 80.0               | 5.75           |                |     | Clear pump, ready  |
|         | 1705 | 110     |          |      | 38-1 1/2       | 84.0               | 7.00           |                |     | resting  |
| 2 hours | 1715 | 120     |          |      | 39-2 1/2       |                    | 7.27           |                |     | Pumped 20 gal  |
|         | 1730 | 135     |          |      | 38-11          |                    | 6.97           |                |     |  |
| 15 min  | 1745 | 150     |          |      | 39-2 1/2       |                    | 7.25           |                |     |  |
|         | 1800 | 165     |          |      | 39-0 1/2       |                    | 7.08           |                |     |  |
| 34      | 1815 | 180     |          |      | 39-2           |                    | 7.21           |                |     | Pumped 30 gal  |
|         | 1830 | 195     |          |      | 39-5           |                    | 7.46           |                |     |  |
|         | 1845 | 210     |          |      | 39-5           |                    | 7.46           |                |     |  |
|         | 1900 | 225     |          |      | 39-2           |                    | 7.21           |                |     |  |
| 44      | 1915 | 240     |          |      | 38-1 1/2       |                    | 7.00           |                |     | Pumped 40 gal  |
|         | 1930 | 255     |          |      | 39-0           |                    | 7.04           |                | 10  |  |
|         | 1932 | 257     |          |      | 39-4           |                    | 7.38           |                | 20  | INCREASE PUMP RATE TO 30 GPM                               |
|         | 1934 | 259     |          |      | 39-10          |                    | 7.87           |                |     | PUMP SLIPPED APPROX 1 FOOT - REPLACED TO APPROX SAME LEVEL |
| 2 min   | 1936 | 261     |          |      | 40-3 1/2       |                    | 8.23           |                |     |  |
|         | 1938 | 263     |          |      | 40-3 1/2       |                    | 8.33           |                |     |  |
|         | 1940 | 265     |          |      | 40-1 1/2       |                    | 8.17           |                |     | PUMPING RATE FALL-OFF SLIGHTLY                             |
|         | 1945 | 270     |          |      | 40-9           |                    | 8.79           |                |     |  |
|         | 1950 | 275     |          |      | 41-9 1/2       |                    | 9.83           |                |     |  |
| 5 min   | 1955 | 280     |          |      | 42-0           |                    | 10.04          |                |     |  |
|         | 2000 | 285     |          |      | 41-8           |                    | 9.72           |                |     |  |
|         | 2005 | 290     |          |      | 42-8 1/2       |                    | 9.75           |                |     |  |
|         | 2010 | 295     |          |      | 43-4           |                    | 11.27          |                |     | Silt in pump   |
| 15 min  | 2015 | 300     |          |      | 43-2 1/2       |                    | 11.25          |                |     | Adj pump valve   |
| 54      | 2030 | 315     |          |      | 43-10 1/2      |                    | 11.92          |                |     | Pumped 10 gal  |

*[Signature]*

# RECOVERY DATA

Page 1 of 1

## DATA SHEET FOR RECORDING PUMP TEST DATA

County: Craig

Observation well no. 24

Location: Craig Refinery

Pumped well no. 24

Average O 51m r = ft. r<sup>2</sup> =

| Date  | Hour  | t (min) | t' (min) | t/t' | Depth to water | s (unadjusted) | Adjustment Δs | s' (adjusted) | Q (gpm) | Remarks |
|-------|-------|---------|----------|------|----------------|----------------|---------------|---------------|---------|---------|
| 2/20  | 20:32 |         | 0        |      | 43-10"         | 144"           | 12.0'         | 0             | 20 gph  |         |
|       | 20:34 |         | 2        |      | 42-7"          |                | 10.63'        | 1.37          |         |         |
|       | 20:36 |         | 4        |      | 41-8 3/4"      |                | 9.69          | 2.31          |         |         |
| 2min  | 20:38 |         | 6        |      | 40-10 1/4"     |                | 8.86          | 3.14          |         |         |
|       | 20:40 |         | 8        |      | 40-2"          |                | 8.21          | 3.79          |         |         |
|       | 20:42 |         | 10       |      | 39-6 1/4"      |                | 7.56          | 4.44          |         |         |
|       | 20:47 |         | 15       |      | 38-0 3/4"      |                | 6.11          | 5.89          |         |         |
|       | 20:52 |         | 20       |      | 37-1 1/2"      |                | 5.17          | 6.83          |         |         |
|       | 20:57 |         | 25       |      | 36'-4"         |                | 4.38          | 7.62          |         |         |
|       | 21:02 |         | 30       |      | 35-8 1/2"      |                | 3.75          | 8.25          |         |         |
|       | 21:07 |         | 35       |      | 35-2 1/4"      |                | 3.27          | 8.73          |         |         |
|       | 21:12 |         | 40       |      | 34-9 1/4"      |                | 2.86          | 9.14          |         |         |
| 5min  | 21:17 |         | 45       |      | 34-5 3/4"      |                | 2.52          | 9.48          |         |         |
|       | 21:22 |         | 50       |      | 34-3 3/4"      |                | 2.27          | 9.73          |         |         |
|       | 21:27 |         | 55       |      | 34-0 1/2"      |                | 2.07          | 9.93          |         |         |
|       | 21:32 |         | 60       |      | 33-10 1/2"     |                | 1.92          | 10.08         |         |         |
| 10min | 21:42 |         | 70       |      | 33-7 1/2"      |                | 1.66          | 10.34         |         |         |
|       | 21:52 |         | 80       |      | 33-5 3/4"      |                | 1.57          | 10.49         |         |         |
|       | 22:02 |         | 90       |      | 33-4 1/2"      |                | 1.42          | 10.58         |         |         |
|       | 22:12 |         | 100      |      | 33-3 3/4"      |                | 1.32          | 10.68         |         |         |
|       | 22:22 |         | 110      |      |                |                |               |               |         |         |
|       | 22:32 |         | 120      |      |                |                |               |               |         |         |
| 15min | 22:47 |         | 135      |      |                |                |               |               |         |         |
|       | 23:02 |         | 150      |      |                |                |               |               |         |         |
|       | 23:17 |         | 165      |      |                |                |               |               |         |         |

TABLE F-2

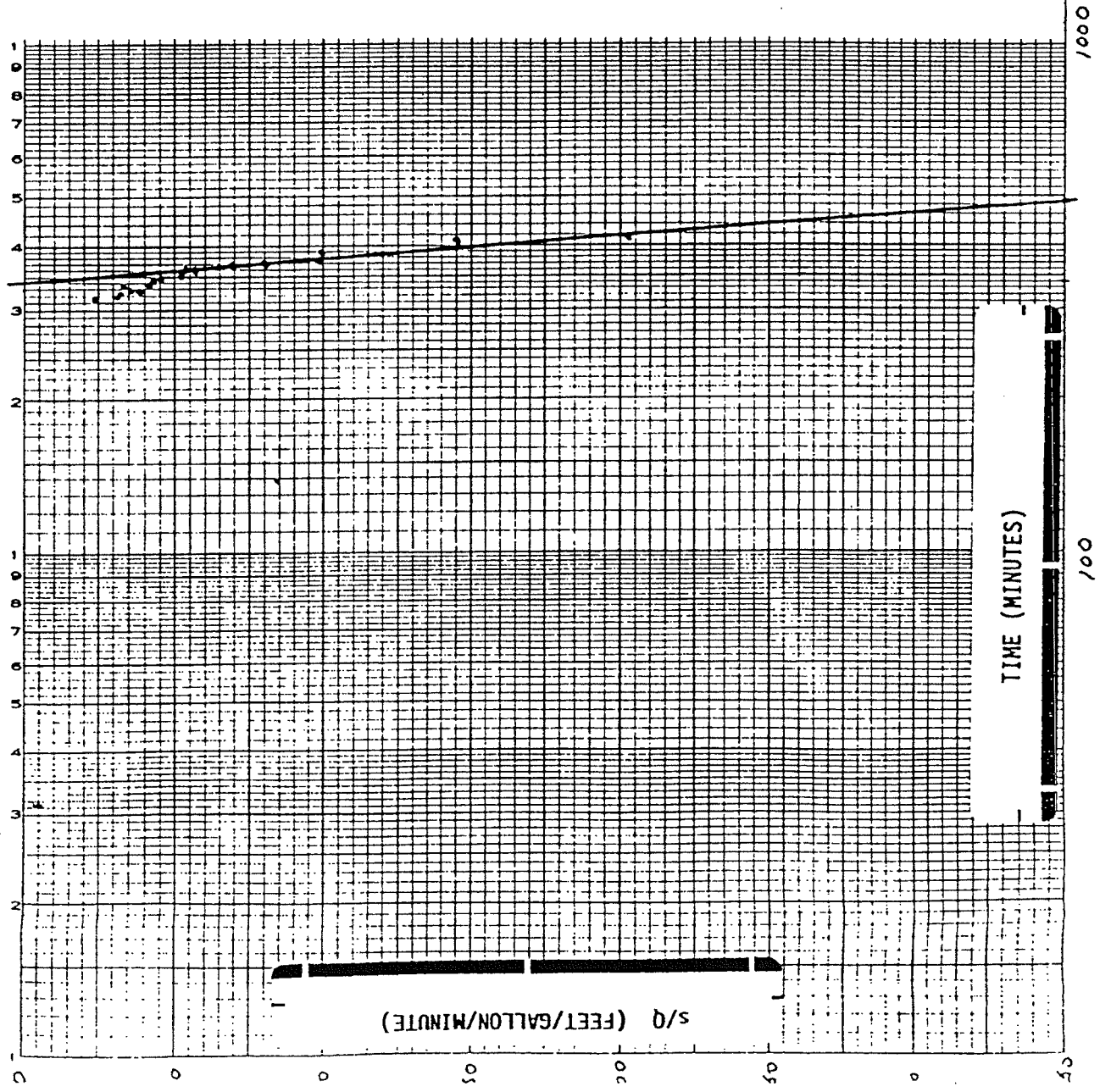
## DATA FOR SHAFER'S METHOD

| Time Since<br>Pumping Started<br>(min) | Drawdown<br>(feet)<br>(s) | Feet of<br>Casing Filled<br>(ft) | Time<br>To Fill<br>(min) | Q<br>(gpm) | S/Q<br>(ft/gpm) |
|--|---------------------------|----------------------------------|--------------------------|------------|-----------------|
| 317                                    | 12.0                      | 0                                | 0                        | ---        | ---             |
| 319                                    | 10.63                     | 1.37                             | 2                        | .45        | 23.8            |
| 321                                    | 9.69                      | .94                              | 2                        | .31        | 31.6            |
| 323                                    | 8.86                      | .83                              | 2                        | .27        | 32.7            |
| 325                                    | 8.21                      | .65                              | 2                        | .21        | 38.7            |
| 327                                    | 7.56                      | .65                              | 2                        | .21        | 35.6            |
| 332                                    | 6.11                      | 1.45                             | 5                        | .19        | 32.3            |
| 337                                    | 5.17                      | .94                              | 5                        | .13        | 42.1            |
| 342                                    | 4.38                      | .79                              | 5                        | .10        | 42.5            |
| 347                                    | 3.75                      | .63                              | 5                        | .08        | 45.6            |
| 352                                    | 3.27                      | .48                              | 5                        | .06        | 52.2            |
| 357                                    | 2.86                      | .41                              | 5                        | .05        | 53.4            |
| 362                                    | 2.52                      | .34                              | 5                        | .04        | 56.8            |
| 367                                    | 2.27                      | .25                              | 5                        | .03        | 69.5            |
| 372                                    | 2.07                      | .20                              | 5                        | .026       | 79.3            |
| 377                                    | 1.92                      | .15                              | 5                        | .019       | 98.0            |
| 387                                    | 1.66                      | .26                              | 10                       | .017       | 97.8            |
| 397                                    | 1.51                      | .15                              | 10                       | .009       | 154             |
| 407                                    | 1.42                      | .15                              | 10                       | .009       | 145             |
| 417                                    | 1.32                      | .10                              | 10                       | .006       | 202             |

FIGURE F-1  
 SHAFER PLOT OF DATA FROM OW-24

Calculation of T & K

- 1) Over the total vertical scale of 350 feet/GPM, time varies by 0.1567 log cycles
- 2) For 1 full log cycle,  $s/Q$  equals  $350/.1567 = 2233$
- 3)  $T = 264/(s/Q) = 264/2233$   
or  $T = 0.110$  gallons/day/foot
- 4) For a 20 foot screened interval  
 $K = T/b = 0.0055$  gallons/day/square foot.
- 5)  $.0055 \text{ g/d/ft} \times 1.55 \times 10^{-6} \text{ (ft/sec)/(g/d/ft)}$   
 $= 8.3 \times 10^{-9} \text{ ft/sec}$



BO. 16 OW-24

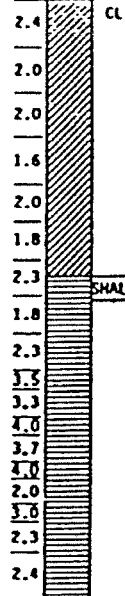
SURFACE ELEVATION: 6676 FEET

| DEPTH IN FEET | LABORATORY TEST DATA     |                  |                      |                    |                                   |                      |                       |                   |
|---------------|--------------------------|------------------|----------------------|--------------------|-----------------------------------|----------------------|-----------------------|-------------------|
|               | TESTS REPORTED ELSEWHERE | ATTENBERG LIMITS |                      | STRENGTH TEST DATA |                                   |                      |                       | DRY DENSITY (PCF) |
|               |                          | LIQUID LIMIT (%) | PLASTICITY INDEX (%) | TYPE OF TEST       | AXIAL OR CONFINING PRESSURE (PSF) | SHEAR STRENGTH (PSF) | DEVIATOR STRESS (PSF) |                   |
| 0             |                          |                  |                      |                    |                                   |                      |                       |                   |
| 10            |                          |                  |                      |                    |                                   |                      |                       |                   |
| 20            |                          |                  |                      |                    |                                   |                      |                       |                   |
| 30            |                          |                  |                      |                    |                                   |                      |                       |                   |
| 40            |                          |                  |                      |                    |                                   |                      |                       |                   |
| 50            |                          |                  |                      |                    |                                   |                      |                       |                   |
| 60            |                          |                  |                      |                    |                                   |                      |                       |                   |
| 70            |                          |                  |                      |                    |                                   |                      |                       |                   |
| 80            |                          |                  |                      |                    |                                   |                      |                       |                   |
| 90            |                          |                  |                      |                    |                                   |                      |                       |                   |
| 100           |                          |                  |                      |                    |                                   |                      |                       |                   |
| 110           |                          |                  |                      |                    |                                   |                      |                       |                   |
| 120           |                          |                  |                      |                    |                                   |                      |                       |                   |
| 130           |                          |                  |                      |                    |                                   |                      |                       |                   |
| 140           |                          |                  |                      |                    |                                   |                      |                       |                   |
| 150           |                          |                  |                      |                    |                                   |                      |                       |                   |
| 160           |                          |                  |                      |                    |                                   |                      |                       |                   |

PENETRATION RATE  
MINUTES/FOOT

SYMBOLS

DESCRIPTION



CL TRIASSIC PERIOD  
CHIMLE FORMATION  
REDDISH BROWN, VERY FINE SANDY CLAY, SOFT, HIGHLY WEATHERED

30 FEET: SHALE, REDDISH BROWN, SANDY, SOFT, FRESH

BORING COMPLETED AT 65.0 FEET ON 1/2/81.  
4-INCH PVC PIEZOMETER INSTALLED WITH PERFORATIONS FROM 41.0 TO 61.0 FEET.  
GRAVEL PLACED FROM 28.0 TO 61.0 FEET AND BORING SEALED WITH BENTONITE AND CEMENT TO SURFACE.  
GROUND WATER LEVEL MEASURED AT 32.5 FEET BELOW GROUND ON 1/5/81.

FIGURE 'F'-2  
LOG OF WELL OW-24

# LOG OF BORINGS

DAMES & MOORE

# Pumping Test Analyses for Low Yield Formations

by David C. Shafer

Occasionally it is necessary to determine aquifer characteristics of very low yielding formations—those with transmissivities less than 500 gallons per day per foot. Though interest in these aquifers is certainly not because of their productive capability, it may be desirable to determine groundwater flow characteristics even in these low yield formations in order to determine such things as regional groundwater flow patterns, effect of dewatering or migration of pollution plumes near point sources of contamination.

## Different Approach

Conventional pumping test analysis using the standard time drawdown graph often does not work effectively in low T (transmissivity) formations for two reasons. First, the pumped well's low specific capacity (gallons per minute per foot of drawdown) may cause the pump to break suction during the test and it may be impractical to throttle back the pumping rate sufficiently to prevent this. Second, even if a constant pumping rate can be maintained without breaking suction, most of the data obtained will probably reflect casing storage effects rather than true aquifer parameters (see "Casing Storage Can Affect Pumping Test Data,"



William F. Achuff  
Director

Jan-Feb. 1978, Johnson Drillers Journal). Thus a different approach is required.

The best method for analyzing these formations is to pump a substantial portion of the casing empty, then shut the pump off and measure water levels as they recover. In ordinary pumping tests these measurements correspond to the non-pumping portion of the test. However, in the low T formations this "recovery period" is actually the "pumping period!"

After pump shut-off, the casing slowly begins filling with water. This water comes from the aquifer and actually represents the water pumped during this so called "pumping period." The pumping rate is determined by measuring the volume of

Pumping rate = 10 gpm  
Pumping period = 15 minutes  
Drawdown at pump shut off = 90 ft  
Casing — 6" I.D.  
Drop pipe — 1¼" I.D.

| Time in minutes<br>since pumping<br>started<br>(t) | Drawdown<br>in feet<br>(s) | Number of<br>feet of<br>casing filled | Time<br>in minutes<br>required<br>to fill | Volume filled<br>divided by time<br>required in gallons<br>per minute<br>(Q) | s/Q<br>in feet<br>per gallon<br>per minute |
|--|----------------------------|---------------------------------------|---|--|--|
| 15<br>(pump shut off)                              | 90                         |                                       |   |  |  |
| 17   | 85.66                      | 4.34                                  | 2   | 3.04   | 28.2                                       |
| 20   | 79.7                       | 5.96                                  | 3   | 2.78   | 28.6                                       |
| 30   | 64.2                       | 15.5                                  | 10  | 2.17   | 29.5                                       |
| 40   | 51.9                       | 12.3                                  | 10  | 1.72   | 30.2                                       |
| 60   | 35.6                       | 13.3                                  | 20  | 1.14   | 31.1                                       |
| 80   | 24.6                       | 11.0                                  | 20  | .77  | 31.8                                       |

Table 1



casing filled in a given length of time.

During the test, careful measurements are made of time since pumping began ( $t$ ) along with drawdown ( $s$ ) at each of these times. Then a calculation is made to determine  $Q$  for each time  $t$  and finally the ratio  $s/Q$  is computed for each measured drawdown value. The ratio is simply the reciprocal of the specific capacity.

A graph is then constructed showing  $t$  versus the ratio  $s/Q$  plotted as usual on semi-logarithmic graph paper with  $t$  on the log scale. A straight line of best fit is drawn through the data points and  $T$  is calculated as follows:

$$T = \frac{264}{\Delta(s/Q)}$$

where  $\Delta(s/Q)$  is the change in  $s/Q$

over one log cycle of graph paper.

This graph has the unique advantage that it will accurately reflect aquifer transmissivity independent of casing storage effects. In addition it will be sensitive to nearby recharge and/or negative boundaries and will reveal these conditions like any ordinary time drawdown graph.

To see how this technique works it is best to work an example. Table 1 shows data obtained from a 6-inch well pumped at 10 gpm for 15 minutes. Drawdown after 15 minutes of pumping measured 90 feet.

The next data point was recorded two minutes following pump shut-off or 17 minutes since pumping started. At this time the pumping water level was 85.66 feet, indicating that 4.34 feet of casing had filled during the two minute interval.

The annulus between the 6-inch casing and 1 1/2" drop pipe holds 1.4 gallons per foot so that the volume of casing filled is 1.4 times 4.34, or 6.08 gallons in two minutes. Thus,

$$Q = 6.08 \text{ gallons}/2 \text{ minutes} = 3.04 \text{ gpm}$$

finally,

$$s/Q = 85.66 \text{ ft}/3.04 \text{ gpm} = 28.2 \text{ ft/gpm}$$

which is plotted at a time of 17 minutes on the graph shown here.

This analysis is repeated for each

drawdown measurement. The resultant calculated  $s/Q$  values are shown in the table and plotted in the figure. The formation  $T$  value from the graph is

$$T = \frac{264}{\Delta(s/Q)} = \frac{264}{5.3} = 49 \text{ gpd/ft}$$

## Conventional Analysis

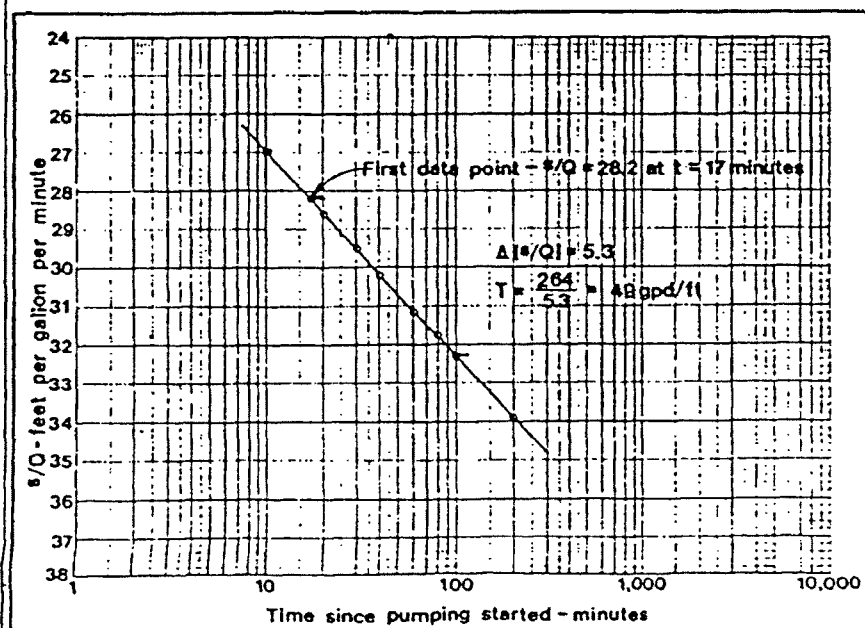
Examination of the hydraulic characteristics of this well (not including the drawdown graph) had been used to show that a conventional time drawdown graph would have lasted for approximately two hours. This means that data recorded in the first twelve hours of pumping would have been useless and longer pumping than this would have been required to obtain any usable data at all. However, data collected after twelve hours of pumping probably would be more influenced by boundary conditions than by aquifer transmissivity. Thus, in practice, perfectly might have been impossible to determine the  $T$  value using conventional analysis techniques regardless of the length of the test. The value of the method described above becomes very clear; it may be the only way to determine  $T$  values in certain low yielding aquifers.

In order to maximize the accuracy of this method, it is best to unload (empty) the casing as rapidly as possible. Thus it is actually better to use a high capacity pump than a low capacity pump in analyzing extremely low-yielding wells!

Another good idea is to unload the casing with compressed air since this can typically be done in one minute or less.

## Recorded Data Must Be Accurate

An additional important consideration is that all data recorded for this type of analysis must be absolutely accurate. Small errors in the recorded values of time and/or drawdown can result in large errors in the calculated values of  $s/Q$ . For best results, drawdown should be recorded to the nearest hundredth of a foot and timed to the nearest second or two.



In low transmissivity situations, readings are taken after pump shut-off. In this method,  $s/Q$  is the reciprocal of the specific capacity and  $t$  is time, measured after shut-off as water begins to enter the casing.



JOHN W. SHOMAKER  
CONSULTING GEOLOGIST  
3236 CANDELARIA, N.E.  
ALBUQUERQUE, NEW MEXICO 87107

(505) 884-2897

September 20, 1984

RECEIVED SEP 24 1984

Carl D. Shook, Plant Manager  
Giant Refining Company, Ciniza Refinery  
Route 3, Box 7  
Gallup, New Mexico 87301

Re: results of permeability tests, July 2 and 3, 1984

Dear Carl:

Copies of the field notes, calculations, and data plots for the two permeability tests are attached. The tests are summarized as follows:

Well OW-4 The well is completed principally in the clay and shale sequence which overlies the uppermost aquifer; a small thickness of sandstone which may be part of the uppermost aquifer was also penetrated. Total depth when drilled was 102.0 ft. Perforations are from 62.0 ft to 102 ft. The well is located near the center of the land-treatment area. A slug test was performed on July 3, 1984, following the method described by S. W. Lohman (1972, Ground-Water Hydraulics, U. S. Geol. Survey Prof. Paper 708, p. 27-29), which indicates the permeability of the section open to the well to be about  $4 \times 10^{-7}$  cm/sec.

Well MW-1 This well is one of the monitoring wells on the boundary of the land-treatment area, and is completed in the uppermost aquifer. It was drilled to 120 ft, and is screened in the interval 87 to 120 ft; the casing is sealed above 89 ft so as to isolate the uppermost aquifer. The slug test performed on July 3, 1984 indicated a permeability of about  $1.2 \times 10^{-4}$  cm/sec.

Information as to the construction of the wells is taken from Dames and Moore (March, 1981; Ground water and soils investigation, Ciniza Refinery near Gallup, New Mexico, and November, 1981, Ground-water monitoring plan, Ciniza Refinery near Gallup, New Mexico).

Please let me know if there are questions.

Sincerely,

John W. Shomaker  
Consulting Geologist

cc: Tom Andrews, Delta H Engineering, W/anal

WCH: JIMMIE - MIZA OW-4  
date: 7-3-84

casing size: nominal 4 1/2" ID 4.0 csg. matl. PVC  
water levels measured from top 4 1/2 OD PVC casing, 5 side,  
which is 1.7 ± above ground level.  
volume of slug:

$$OD, ft \left( \frac{0.27}{0.30} \times \frac{1}{2} \right)^2 \times \pi \times \text{length, ft} \frac{0.10}{0.04} = \frac{0.006}{0.003} ft^3$$

|              |             |              |
|--------------|-------------|--------------|
| <u>0.243</u> | <u>4.84</u> | <u>0.224</u> |
| <u>0.27</u>  | <u>0.26</u> | <u>0.015</u> |
| <u>0.243</u> | <u>4.81</u> | <u>0.223</u> |
| <u>0.27</u>  | <u>0.26</u> | <u>0.015</u> |
| <u>0.18</u>  | <u>0.03</u> | <u>0.001</u> |
| <u>0.11</u>  | <u>0.25</u> | <u>0.002</u> |
|              |             | <u>/</u>     |

Note - obstruction  
in casing at 33'  
won't pass 0.27  
OD.

Sum 0.489 =  $V, ft^3 = 3.66 \text{ gal.}$

$r_c$  = internal radius of casing above perms: 0.165 ft

$r_s$  = radius of screen or open hole: \_\_\_\_\_ ft

initial water level 26.15 ft below MP, time 07:51

$$H_0 = \frac{V}{\pi r_c^2} = \frac{0.489}{\pi (0.165)^2} = \underline{5.72} \text{ ft}$$

| clock time | t, sec. | water level | H    | H/H <sub>0</sub> | remarks             |
|------------|---------|-------------|------|------------------|---------------------|
|            | 0       |             |      |                  |                     |
| 08:20:54   |         |             |      |                  | slug released       |
| 21:25?     | 31      | 20.75       | 5.40 | 0.944            | top slug: 23.2' MP  |
| 22:28      | 94      | 20.76       | 5.39 | 0.942            |                     |
| 23:14      | 140     | 20.80       | 5.35 | 0.935            |                     |
| 23:51      | 177     | 20.81       | 5.34 | 0.934            |                     |
| 24:42      | 228     | 20.83       | 5.32 | 0.930            |                     |
| 25:42      | 288     | 20.87       | 5.28 | 0.923            |                     |
| 27:06      | 372     | 20.90       | 5.25 | 0.918            |                     |
| 28:14      | 440     | 20.92       | 5.23 | 0.914            |                     |
| 29:43      | 529     | 20.95       | 5.20 | 0.909            |                     |
| 30:55      | 601     | 20.97       | 5.18 | 0.906            |                     |
| 32:46      | 712     | 21.00       | 5.15 | 0.900            |                     |
| 35:16      | 862     | 21.04       | 5.11 | 0.893            |                     |
| 38:38      | 1064    | 21.10       | 5.05 | 0.883            |                     |
| 43:20      | 1356    | 21.19       | 4.96 | 0.867            |                     |
| 46:52      | 1558    | 21.21       | 4.94 | 0.864            | raised slug 0.4 ft. |
| 49:26      | 1712    | 21.25       | 4.90 | 0.857            |                     |
| 52:54      | 1920    | 21.30       | 4.85 | 0.848            |                     |



OLETZG ENCORPORATION

ow-4  
7-3-84

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# maths

$$\frac{W}{W_0} = 1$$

date: 7-2-84

casing size: nominal  $5\frac{1}{2}$ " OD, ID 5.1" csg. matl. PVC  
 water levels measured from top PVC casing, SW side  
 which is 1.37 ft above ground level, (concrete slab)  
 volume of slug: 2.05

$$\text{cap OD, ft} \left( \frac{0.37}{2} \right)^2 \times \pi \times \text{length, ft} = \text{ft}^3$$

|      |      |       |
|------|------|-------|
| 0.39 | 0.05 | 0.006 |
| 0.33 | 4.81 | 0.411 |
| 0.38 | 0.28 | 0.032 |
| 0.33 | 0.04 | 0.003 |
| 0.42 | 0.04 | 0.006 |
| 0.29 | 0.03 | 0.002 |
| 0.16 | 0.03 | 0.001 |
| 0.11 | 0.25 | 0.002 |

sum  $0.474 = V, \text{ ft}^3 = 3.55 \text{ gal}$

$r_c$  = internal radius of casing above perms. 0.211 ft

$r_s$  = radius of screen or open hole: \_\_\_\_\_ ft

initial water level 5.72 ft below mp, time 13:32

$$H_0 = \frac{V}{\pi r_c^2} = \frac{0.474}{\pi (0.211)^2} = 3.389 \text{ ft}$$

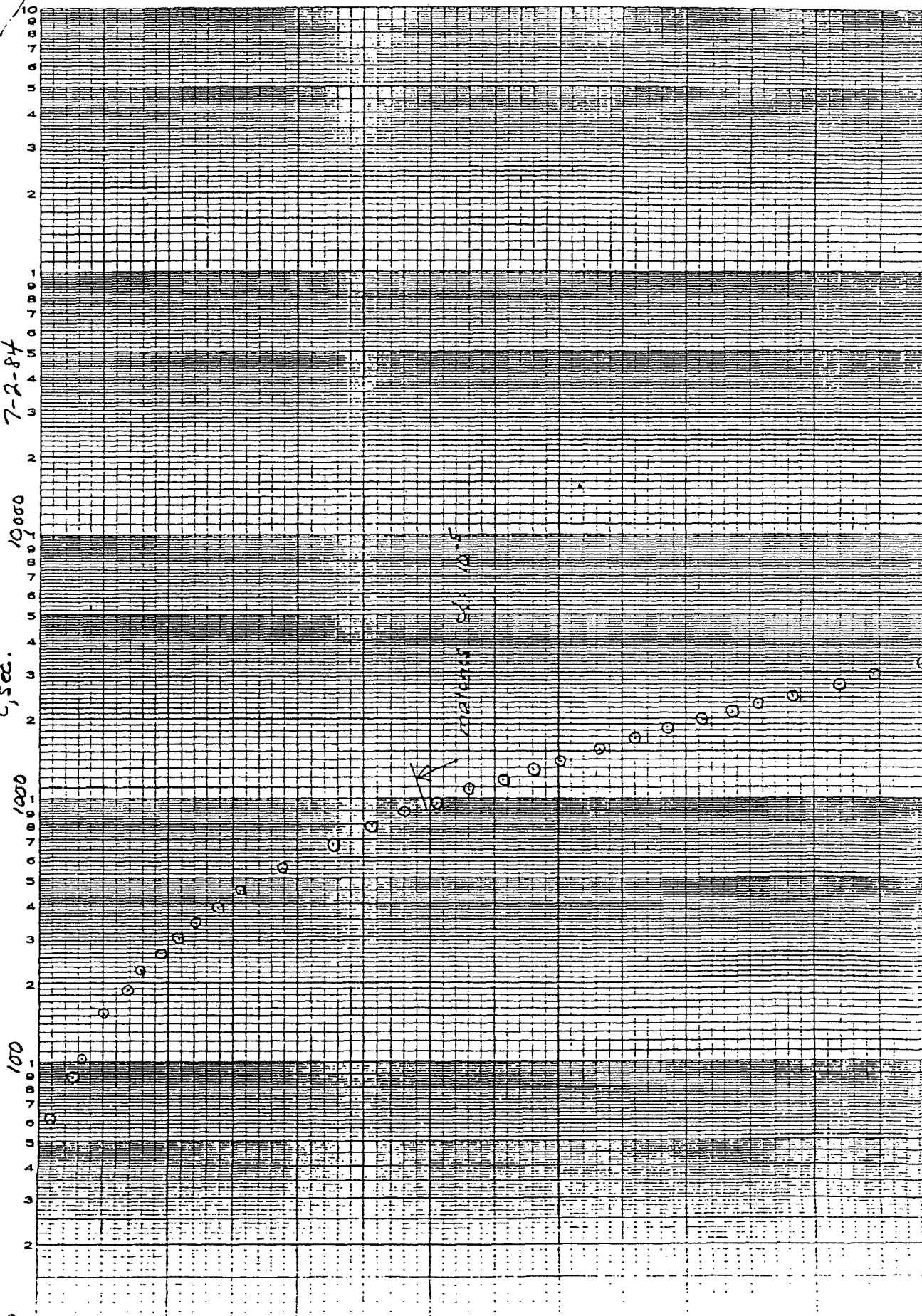
| clock time | $t, \text{ sec.}$ | water level | H    | $H/H_0$ | remarks       |
|------------|-------------------|-------------|------|---------|---------------|
| 13:48:00   | 0                 |             |      |         | slug released |
| 13:48:20   | 20                | 2.30 ft     | 3.42 | 1.009   | ??            |
| 49:00      | 60                | 2.36        | 3.36 | 0.991   |               |
| 49:20      | 80                | 2.42        | 3.30 | 0.974   |               |
| 49:52      | 112               | 2.44        | 3.28 | 0.968   |               |
| 50:32      | 152               | 2.50        | 3.22 | 0.950   |               |
| 51:05      | 185               | 2.56        | 3.16 | 0.932   |               |
| 51:42      | 222               | 2.60        | 3.12 | 0.921   |               |
| 52:19      | 259               | 2.65        | 3.07 | 0.906   |               |
| 52:55      | 295               | 2.70        | 3.02 | 0.891   |               |
| 53:33      | 333               | 2.74        | 2.98 | 0.879   |               |
| 54:20      | 386               | 2.80        | 2.92 | 0.862   |               |
| 55:30      | 450               | 2.86        | 2.86 | 0.844   |               |
| 57:05      | 545               | 2.97        | 2.75 | 0.811   |               |
| 59:12      | 672               | 3.10        | 2.62 | 0.773   |               |
| 14:01:05   | 785               | 3.20        | 2.52 | 0.744   |               |
| 02:25      | 868               | 3.28        | 2.44 | 0.720   |               |
| 03:52      | 952               | 3.36        | 2.36 | 0.696   |               |

[illegible]

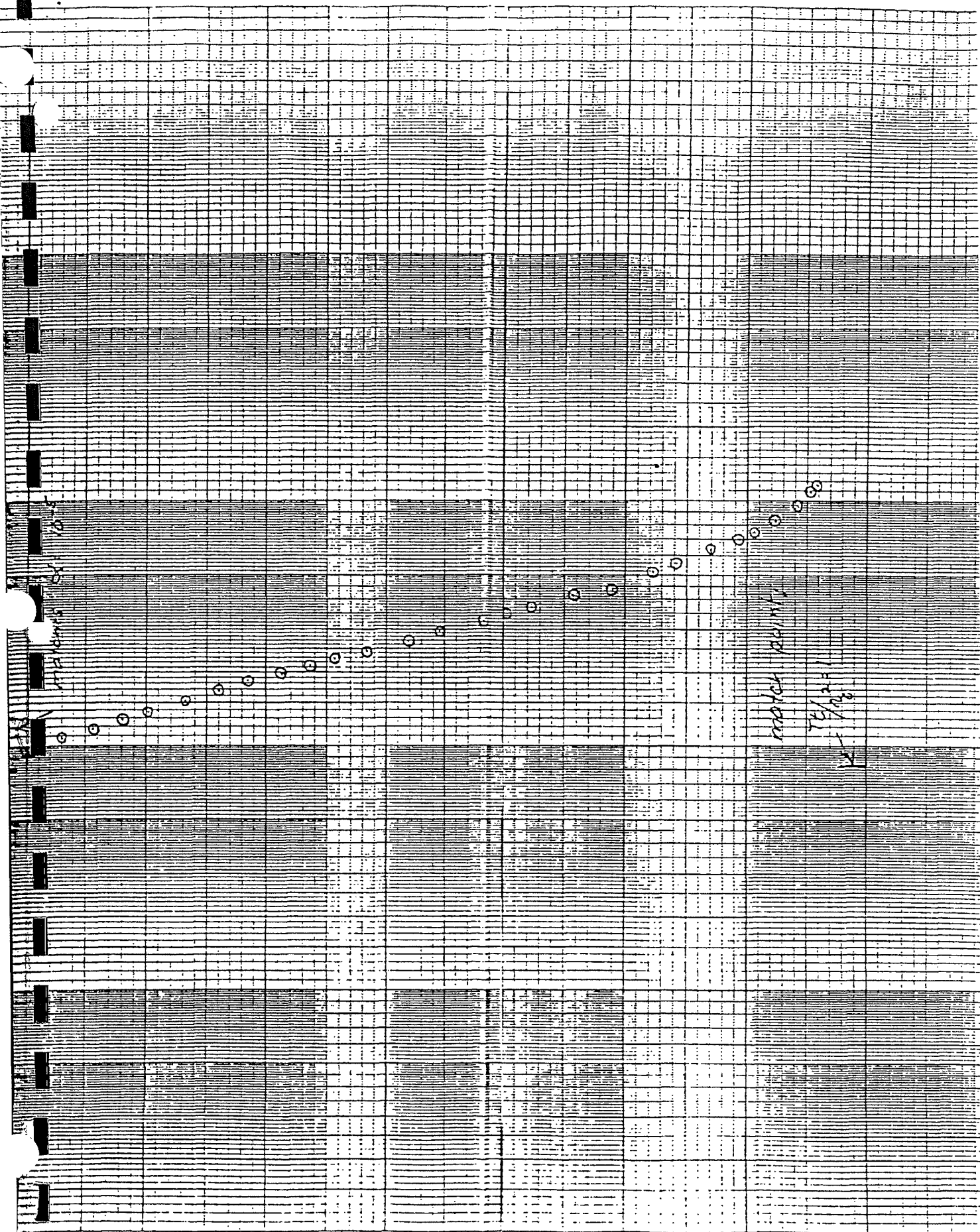
DIETZEN CORPORATION  
MADE IN U.S.A.

4,500.

mw-1  
7-2-84









APPENDIX C  
RCRA POST-CLOSURE CARE PERMIT APPLICATION  
LAND TREATMENT UNIT  
GIANT REFINING COMPANY  
CINIZA REFINERY

May 2000

*RCRA Part A and Part B  
Post-Closure Permit Application  
Land Treatment Unit  
May 2000*

**APPENDIX C**

**Land Treatment Unit Historical Information and Data**

## LAND TREATMENT UNIT HISTORICAL INFORMATION AND DATA

### 1.0 LAND TREATMENT HISTORY

Historical LTU information and data extracted from existing permit applications, operating permits, operating records, and other source documents are provided as Appendix C. The inclusion of this appendix does not imply that historical information and data have been verified.

In August 1980, Ciniza Refinery (Ciniza) notified the U.S. Environmental Protection Agency (EPA) that it was a generator and operator of a hazardous waste management facility. In November 1980, Ciniza submitted a Part A permit application as an "existing facility" (defined at 40 *Code of Federal Regulations* §260.10). This granted Ciniza interim status for their Land Treatment Unit (LTU) operations. In response to notice from the Regional Administrator, Ciniza submitted a Part B permit application in December 1983. Based on changing guidance, Ciniza submitted a land treatment demonstration plan (LTD) and an application for a two-phase LTD permit in April 1985. On February 9, 1987, Ciniza was issued a Short-term LTD Permit (NMD000333211-1) to conduct a hazardous waste land LTD. The LTD was conducted to identify the land treatment capabilities for refinery waste generated by Ciniza. The LTD defined waste management parameters (e.g., rate-limiting constituent, application-limiting constituent, capacity-limiting constituent, and unit life of the LTU). This was accomplished by identifying the Principal Hazardous Constituents (PHCs) present in refinery waste streams and measuring their degradation, transformation, and immobilization in the treatment zone of the LTU. From the results of the LTD and a modified Part B permit application, Ciniza was issued a Hazardous Waste Facility Permit (NMD 000333211-2) on November 4, 1988. Ciniza has not applied hazardous wastes to the LTU since November 8, 1990.

#### 1.1 Land Treatment Program [20 NMAC 4.1, Subpart IX, §270.20(b)]

Ciniza's Hazardous Waste Facility Permit established operational requirements for the LTU. These requirements include procedural and engineering controls necessary to ensure that hazardous constituents are fully treated within the LTU without uncontrolled release to the environment.

The LTU consists of a treatment zone of soil extending 5 ft deep from the original soil surface. This depth is shallow enough to ensure that the treatment zone is more than 3 ft above the seasonal high water table. The zone of incorporation (ZOI) within the treatment zone is the volume of soil to which the waste was directly applied.

The ZOI for the Ciniza LTU is the top 12 in. of the treatment zone. The LTU was designed and constructed to prevent both washout of any hazardous waste and to prevent inundation of and discharge from the permitted unit through the use of a continuous dike which surrounds the LTU at an elevation of 3 ft above the natural grade.

The ZOI was tilled during permitted operations to encourage aerobic microbial activity and improve chemical reaction rates. During active treatment soil nutrients were applied, as necessary, to optimize carbon:nitrogen:phosphorous (C:N:P) ratios. Applications of Ciniza wastes to the LTU were limited to ensure that treatment processes were not overwhelmed or poisoned. Performance indicators (e.g., soil moisture, pH, total organic carbon) were monitored in the ZOI to ensure that treatment was proceeding.

#### **1.2 Treatment Zone Description [20 NMAC 4.1, Subpart IX, §270.20(b)(2) and §270.20(b)(5)]**

The LTU consists of three 480-by-240-ft sections, each of which contain 2.6 acres (1.0 hectares) of available treatment surface. Each section is delineated by a continuous dike to prevent site runoff and runoff. The treatment zone extends 5-ft deep from the top of the soil within the diked section. The top 12 in. of the treatment zone is the ZOI. The ZOI is tilled when active to encourage aerobic degradation of organics and to maintain moisture content of the soil. This leaves 4 ft of the treatment zone undisturbed.

The soil within the treatment zone is silty clay containing closely-spaced root systems in the uppermost 3 to 4 ft. Field infiltration rates (the rate at which water penetrates into the soil surface) averages  $1.0 \times 10^{-3}$  cm/sec or 3.6 cm/hr. Soil permeability as determined by laboratory measurements averages  $1.9 \times 10^{-5}$  cm/sec or  $6.8 \times 10^{-2}$  cm/hr for three locations at the 6- to 12-in. depth. Field infiltration rate allows prediction of runoff and erosion; permeability (hydraulic conductivity) allows estimation of vertical water movement rates in the soil. The treatment zone soils have a saturated hydraulic conductivity rating of "moderately low" by the U.S. Department of Agriculture Class (Giant Refining Company Part B Permit Application 1984). The low permeability of the treatment zone soil assists in retarding the vertical movement of hazardous constituents through the treatment zone.

The silty clay soil has a high cation exchange capacity (CEC). The CEC is the total amount of exchangeable cations that the soil has to exchange with cations in the soil solution. The exchangeable cations in the LTU are the heavy metals present in the Ciniza wastes. The high CEC results in high sorption of heavy metals in the LTU soils, assuming other factors (such as soil pH) are favorable.

RECENT SOIL PERMEABILITY TESTS  
FIREWATER POND CONSTRUCTION PROJECT

Precision Engineering, Inc.

P.O. Box 422  
Las Cruces, NM 88004  
505-523-7674

Rigid Wall Hydraulic Conductivity  
Falling Head

GRCC

Proposed Fire Water  
Storage

ATTN: James Romero  
Giant Refining Company  
Route 3, Box 7  
Gallup, NM 87301

Project: Ciniza Fire Water Lagoon

File No.: 05-100

Soil Type: Silty Clay

Date: October 13, 2005

Lab No.: 47872

Sampled From: Boring 05-100-1(2.5'-3.0')

Performed By: GG

TEST SPECIMEN CONDITIONS AT BEGINING OF TEST:

Wet Unit Weight: 120.8 pcf

% Moisture: 10.8

Dry Unit Weight: 109.0 pcf

% Compaction: n/a

% Compaction Requested: n/a

PROCTOR INFORMATION:

Maximum Dry Density: n/a pcf

Optimum Moisture Content: n/a %

Coefficient of Permeability,  $k_{20}$ :  $1.1 \times 10^{-7}$  cm/sec. avg

Remarks: Avg of three:  $1.1 \times 10^{-7}$ ,  $1.1 \times 10^{-7}$ ,  $1.1 \times 10^{-7}$

C:\bill\Projects\2005\05100cinizafirewtr\Permeability.xls\Report

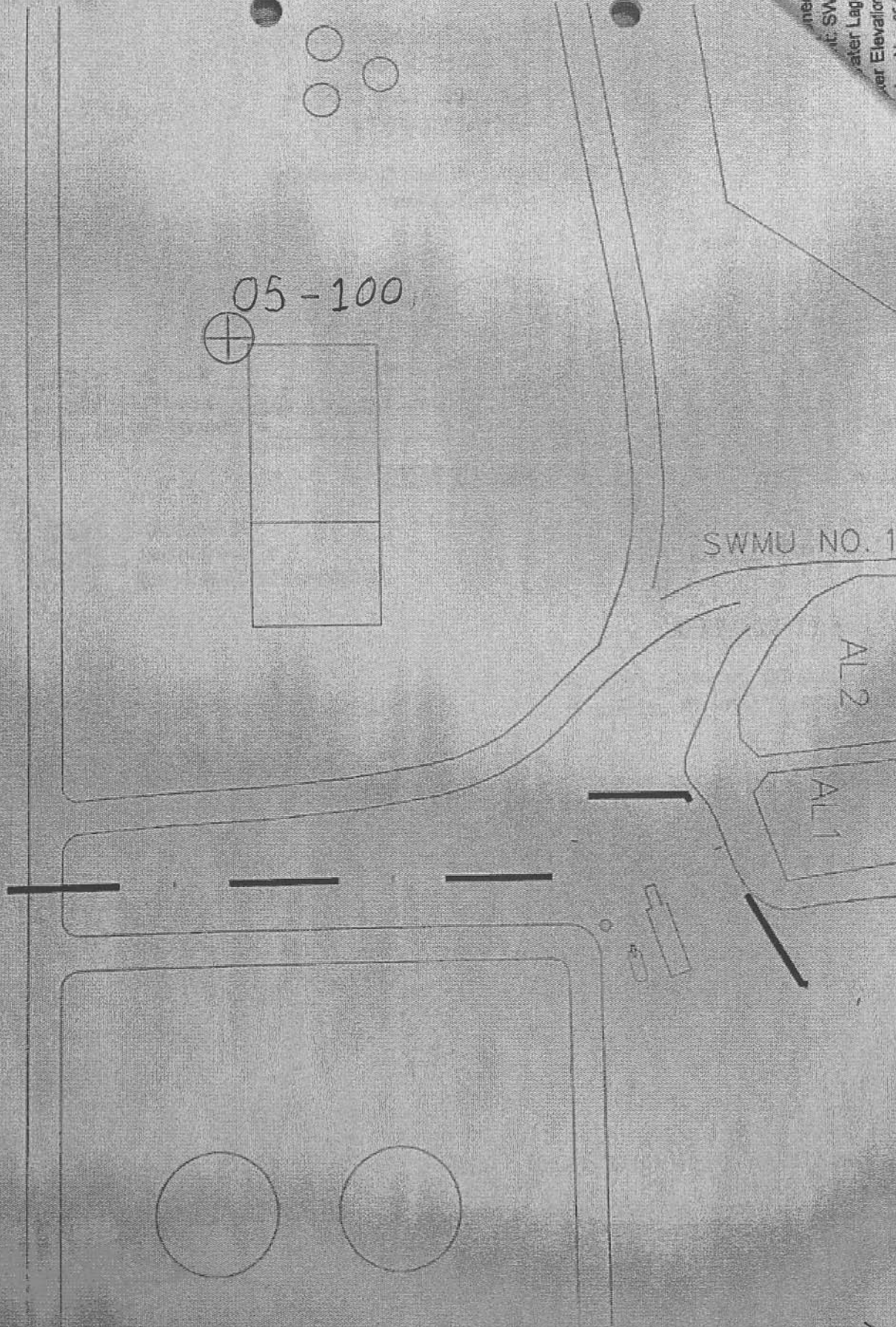
Reviewed By:

Reviewed By:

Certified By:







Sheet: 107  
Site: SW Con  
Water Lagoon  
Water Elevation: 45.15  
Soring No.: 05-100-1

| BLOW  |         |
|-------|---------|
| LAB # | DEPTH   |
| COUN  | 0.0-5.0 |



SW Corner of Proposed  
Lagoon  
Elevation: -88.2@TD, -27@1 hr  
No. 05-100-1

Precision Engineering, Inc.  
P.O. Box 422  
Las Cruces, NM 88004  
505-523-7674

File #: 05-100  
Site: Giant-Ciniza

Elevation: EXISTING  
Date: 9/24/2005

# Log of Test Borings

| LAB # | DEPTH     | BLOW<br>COUNT | PLOT | SCALE | MATERIAL CHARACTERISTICS<br>(MOISTURE, CONDITION, COLOR, ETC.)                                       | %M | LL | PI | CLASS |
|-------|-----------|---------------|------|-------|--|----|----|----|-------|
|       |           |               |      |       |  |    |    |    |       |
| 47872 | 0.0-5.0   |               |      |       | Clay, Very Silty, Sandy, Very Fine, Dark Red,<br>Wet, Firm   |    |    |    |       |
|       |           |               |      | 2.5   | [2.5-3.0 hydraulic conductivity sample]  |    |    |    |       |
|       | 5.0-10.0  |               |      | 5.0   | Same As Above, Wet, Soft   |    |    |    |       |
|       |           |               |      | 7.5   |  |    |    |    |       |
|       | 10.0-15.0 |               |      | 10.0  | Same As Above  |    |    |    |       |
|       | 15.0-16.0 |               |      | 15.0  | Same As Above  |    |    |    |       |
|       | 16.0-17.5 |               |      |       | Sand, Very Fine, Very Clayey, Very Silty,<br>Weak Water Bearing, Moderately Dense, Dark<br>Red/Brown |    |    |    |       |
|       | 17.5-21.5 |               |      | 20.0  | Clay, Dark Red, Wet, Soft  |    |    |    |       |

SIZE & TYPE OF BORING: 4 1/4" ID HOLLOW STEMMED AUGER

LOGGED BY: WHK

C:\bill\Projects\2005\05100\ciniza\firewtr\Boring 1.xls\Sheet1



Sheet: 2 OF 5  
 Bore Point: SW Corner of Proposed  
 Fire Water Lagoon  
 Water Elevation:  
 Boring No.: 05-100-1

Precision Engineering, Inc.  
 P.O. Box 422  
 Las Cruces, NM 88004  
 505-523-7674

File # \_\_\_\_\_  
 Site: SW Corner of Proposed  
 Fire Water Lagoon  
 Elevation: EXISTING  
 Date: 9/24/2005  
 Boring No.: 05-100-1

Log of Test Borings

| LAB # | DEPTH     | BLOW<br>COUNT | PLOT | SCALE | MATERIAL CHARACTERISTICS<br>(MOISTURE, CONDITION, COLOR, ETC.)  | %M | LL | PI | CLASS |
|-------|-----------|---------------|------|-------|---|----|----|----|-------|
|       | 21.5-22.1 |               |      |       | Clay, Sandy, Dark Brown, Wet, Stiff   |    |    |    |       |
|       | 22.1-23.5 |               |      |       | Clay, Hard, Red/Brown (Brighter than Above),<br>Wet/Moist   |    |    |    |       |
|       | 23.5-25.1 |               |      |       | Pertrified Forest Formation<br>Mudstone/Claystone, Weathered, Some<br>Grey/Green Reduction Spots, Generally Red/<br>Brown, Fissile to Crumbly, Damp/Moist |    |    |    |       |
|       | 25.0-30.0 |               |      | 25.0  | Mudstone, As Above, Few Reduction Spots,<br>Damp/Dry  |    |    |    |       |
|       | 30.0-35.0 |               |      | 30.0  | Same As Above dry   |    |    |    |       |
|       | 35.0-40.0 |               |      | 35.0  | Same As Above<br>dry  |    |    |    |       |
|       | 40.0-45.0 |               |      | 40.0  | Same As Above, Brighter Red @ 44.5'-45.0',<br>dry   |    |    |    |       |
|       | 41.5      |               |      |       | siltstone/mudstone, dry, very dense, bright<br>red brown  |    |    |    |       |

SIZE & TYPE OF BORING: 4 1/4" ID HOLLOW STEMMED AUGER  
 C:\blm\Projects\2005\05100\ch\zafirewtr\Boring 1.xls]Sheet1A

LOGGED BY: WHK



Sheet 2 OF 5  
 Bore Point: SW Corner of Proposed  
 Fire Water Lagoon  
 Water Elevation:  
 Boring No.: 05-100-1

Precision Engineering, Inc.  
 P.O. Box 422  
 Las Cruces, NM 88004  
 505-523-7674

File # meet: 3  
 Site: SW Cor  
 Water Lagoon  
 Elevation: EXISTING  
 Date: 9/24/2005  
 Boring No.: 05-100-

### Log of Test Borings

| LAB #   | DEPTH     | BLOW<br>COUNT | PLOT | SCALE | MATERIAL CHARACTERISTICS<br>(MOISTURE, CONDITION, COLOR, ETC.)   | %M | LL | PI | CLASS |
|---|-----------|---------------|------|-------|--|----|----|----|-------|
|   | 21.5-22.1 |               |      |       | Clay, Sandy, Dark Brown, Wet, Stiff  |    |    |    |       |
|   | 22.1-23.5 |               |      |       | Clay, Hard, Red/Brown (Brighter than Above),<br>Wet/Moist  |    |    |    |       |
|   | 23.5-25.1 |               |      |       | Peritrified Forest Formation<br>Mudstone/Claystone, Weathered, Some<br>Grey/Green Reduction Spots, Generally Red/<br>Brown, Fissile to Crumbly, Damp/Moist |    |    |    |       |
|   | 25.0-30.0 |               |      | 25.0  | Mudstone, As Above, Few Reduction Spots,<br>Damp/Dry   |    |    |    |       |
|   | 30.0-35.0 |               |      | 30.0  | Same As Above dry  |    |    |    |       |
|   | 35.0-40.0 |               |      | 35.0  | Same As Above<br>dry   |    |    |    |       |
|   | 40.0-45.0 |               |      | 40.0  | Same As Above, Brighter Red @ 44.5'-45.0',<br>dry  |    |    |    |       |
|   | 41.5      |               |      |       | siltstone/mudstone, dry, very dense, bright<br>red brown   |    |    |    |       |
| SIZE & TYPE OF BORING: 4 1/4" ID HOLLOW STEMMED AUGER |           |               |      |       |  |    |    |    |       |
| LOGGED BY: WHK  |           |               |      |       |  |    |    |    |       |



3 OF 5  
 Corner of Proposed  
 Elevation  
 No. 05-100-1

Precision Engineering, Inc.  
 P.O. Box 422  
 Las Cruces, NM 88004  
 505-623-7874

File #: 05-100  
 Site: Giant Cinza

Elevation: EXISTING  
 Date: 9/24/2005

Log of Test Borings

| LAB #   | DEPTH     | BLOW<br>COUNT | PLOT | SCALE | MATERIAL CHARACTERISTICS<br>(MOISTURE, CONDITION, COLOR ETC.) | %M | LL | PI | CLASS. |
|---|-----------|---------------|------|-------|---|----|----|----|--------|
|   | 45.0-50.0 |               |      | 45.0  | Mudstone/Siltstone, Bright Red/Brown, Dry<br>Very Dense       |    |    |    |        |
|   | 50.0-55.0 |               |      | 50.0  | Same As Above   |    |    |    |        |
|   | 55.0-60.0 |               |      | 55.0  | Same As Above   |    |    |    |        |
|   | 60.0-65.0 |               |      | 60.0  | Same As Above   |    |    |    |        |
| SIZE & TYPE OF BORING: 4 1/4" ID HOLLOW STEMMED AUGER |           |               |      |       |   |    |    |    |        |
| LOGGED BY: WHK  |           |               |      |       |   |    |    |    |        |



Sheet: 4 OF 5  
 Bore Point: SW Corner of Proposed  
 Fire Water Lagoon  
 Water Elevation:  
 Boring No.: 05-100-1

Precision Engineering, Inc.  
 P.O. Box 422  
 Las Cruces, NM 88004  
 505-523-7674

File: 05-100-1  
 Sheet: 4 of 5  
 Site: SW Corner of  
 Fire Water Lagoon  
 Elevation: EXISTING  
 Date: 9/24/2005  
 Boring No.: 05-100-1

Log of Test Borings

| LAB # | DEPTH     | BLOW<br>COUNT | PLOT | SCALE | MATERIAL CHARACTERISTICS<br>(MOISTURE, CONDITION, COLOR, ETC.)  | %M | LL | PI | CLASS |
|-------|-----------|---------------|------|-------|---|----|----|----|-------|
|       | 65.0-66.2 |               |      | 65.0  | Same As Above   |    |    |    |       |
|       | 66.2-67.4 |               |      |       | Sandstone, Fine, Red/Brown, Hard, Fissile,<br>Grey/Green Orbicular Reduction Spots<br>Abundant, Dry                 |    |    |    |       |
|       | 67.4-78.0 |               |      | 70.0  | Sandstone/Siltstone Interbeds, Some<br>Mudstone, Dark Red/Brown, Hard, Dry  |    |    |    |       |
|       |           |               |      | 75.0  |   |    |    |    |       |
|       | 78.0-79.1 |               |      |       | Mudstone, Fissile, Dry, Red/Brown, Some<br>Grey/Green Reduction Spots   |    |    |    |       |
|       | 79.1-97.7 |               |      | 80.0  | Sandstone, Fine, Red/Brown, Fissile, Hard,<br>Feldspathic, Bedded 6" to 1.0', White Mottled<br>Upper Sonsela Member |    |    |    |       |
|       |           |               |      | 85.0  |   |    |    |    |       |

SIZE & TYPE OF BORING: 4 1/4" ID HOLLOW STEMMED AUGER

LOGGED BY: WHK



TICS  
 (RETC)  
 %M  
 LL  
 PI  
 CLASS

Elevation: EXISTING  
 Date: 9/24/2005  
 File # 05-100-1  
 Site: Lagoon  
 Corner of Proposed

Precision Engineering, Inc.  
 P.O. Box 422  
 Las Cruces, NM 88004  
 505-523-7674

File #: 05-100  
 Site: Giant-Ciniza  
 Elevation: EXISTING  
 Date: 9/24/2005

Log of Test Borings

| LAB # | DEPTH      | BLOW COUNT | PLOT | SCALE | MATERIAL CHARACTERISTICS (MOISTURE, CONDITION, COLOR, ETC.)   | %M | LL | PI | CLASS |
|-------|------------|------------|------|-------|---|----|----|----|-------|
|       |            |            |      | 90.0  | Upper Sonsele Member Continued (dry)  |    |    |    |       |
|       |            |            |      | 95.0  |   |    |    |    |       |
|       | 97.7-98.2  |            |      |       | <u>Mudstone</u> , Hard, Dry, Green/Grey-White   |    |    |    |       |
|       | 98.2-101.6 |            |      | 100.0 | <u>Sandstone</u> , Fine to Medium, Quartz Grains, Water Bearing, Hard, (Sonsele Member, Petrified Forest Formation, Chinle Group)   |    |    |    |       |
|       | 101.6      |            |      | 105.0 | TD<br>Boring continuously sampled using 5' split barreled intrusion sampler<br>Boring closed using 10' of 3/8" TR-30 Pel Plug capped with 50' of 8% bentonite cement slurry and backfilled to the ground surface with cuttings. |    |    |    |       |

SIZE & TYPE OF BORING: 4 1/4" ID HOLLOW STEMMED AUGER

LOGGED BY: WHK