

**GW - \_\_\_\_\_1\_\_\_\_\_**

# **WORK PLANS**

**Investigation Work  
Plan Group 6  
12/2009**

October 9, 2012

2012 OCT 12 AM 10:55

John E. Kieling, Chief  
New Mexico Environment Department  
Hazardous Waste Bureau  
2905 Rodeo Park Drive East, Bldg 1  
Santa Fe, NM 87505

**Certified Mail #: 7009 0820 0000 0482 9238 (Delivery to NMED)****Certified Mail #: 7010 1870 0002 6760 0139 (Delivery to OCD)**

RE: Response to NMED's Approval  
Investigation Report Group 6 (AOC No. 19 Seep North of MW-45, AOC No. 20 Seep North of MW-46, and AOC No. 21 Seep North of MW-47)  
Western Refining Southwest, Inc., Bloomfield Refinery  
EPA ID# NMD089416416

Dear Mr. Kieling:

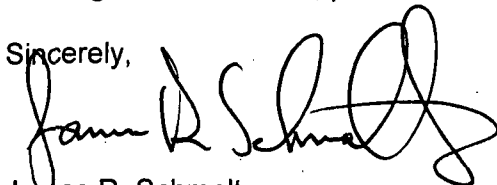
Pursuant to your approval letter dated July 5, 2012 for the Group 6 Investigation Report, Western Refining Southwest, Inc. - Bloomfield Refinery ("Western") has prepared the following letter Investigation Work Plan in response to your comment No. 2 (shown below).

**NMED Comment 2:**

*Section 7.2 (Recommendations), page 34, states that "[a]dditional soil samples should be collected at the locations with the highest concentrations of chromium at AOC 19 [e.g., AOC 19-13 (0-0.5'), AOC 19-14 (0-0.5'), and AOC 19-18 (0-0.5')]. The samples should be analyzed for chromium VI to determine if the chromium detected in soils at AOC 19 is present as chromium III or chromium VI." Western must collect at least three additional soil samples from soils within AOC 19 at locations where contaminants are not detected at high concentrations to determine if chromium is typically present as chromium III and not present only as a result of local reducing conditions in areas containing petroleum contamination. Western must submit a letter work plan to propose the additional investigation activities to collect these soil samples.*

If you have questions regarding the proposed investigation presented in the attached Investigation Work Plan, please contact me at (505) 632-4171.

Sincerely,



James R. Schmaltz  
Health, Safety, Environmental, and Regulatory Director  
Western Refining Southwest, Inc., Bloomfield Refinery

Enclosure

cc: Dave Cobrain – NMED HWB  
Leona Tsinnajinnie – NMED HWB

✓ Carl Chavez - NMOCD  
Allen Hains – Western Refining El Paso

# **GROUP 6 (AOC No. 19 – SEEP NORTH OF MW-45) INVESTIGATION WORK PLAN ADDENDUM**

## **Introduction**

The Bloomfield Refinery is located immediately south of Bloomfield, New Mexico in San Juan County. The physical address is #50 Road 4990, Bloomfield, New Mexico 87413. The Bloomfield Refinery is located on approximately 263 acres. Bordering the facility is a combination of federal and private properties. Public property managed by the Bureau of Land Management lies to the south. The majority of undeveloped land in the vicinity of the facility is used extensively for oil and gas production and, in some instances, grazing. U.S. Highway 550 is located approximately one-half mile west of the facility. The topography of the main portion of the site is generally flat with steep bluffs to the north where the San Juan River intersects Tertiary terrace deposits.

The Bloomfield Refinery is a crude oil refinery currently owned by Western Refining Southwest, Inc., which is a wholly owned subsidiary of Western Refining Company, and it is operated by Western Refining Southwest, Inc. – Bloomfield Refinery. The Bloomfield Refinery has an approximate refining capacity of 18,000 barrels per day. Various process units operated at the facility, including crude distillation, reforming, fluidized catalytic cracking, sulfur recovery, merox treater, catalytic polymerization, and diesel hydrotreating. Past operations produced gasoline, diesel fuels, jet fuels, kerosene, propane, butane, naphtha, residual fuel, fuel oils, and LPG.

On July 27, 2007, the New Mexico Environment Department (NMED) issued an Order to San Juan Refining Company and Giant Industries Arizona, Inc. ("Western") requiring investigation and corrective action at the Bloomfield Refinery. This Investigation Work Plan has been prepared for the Areas of Concern (AOC) designated as Group 6 in the Order. Group 6 includes AOC No. 19 (Seep North of MW-45), AOC No. 20 (Seep North of MW-46), and AOC No. 21 (Seep North of MW-47). The location of AOC Nos. 19, 20 and 21 are shown on Figure 1. This Investigation Work Plan focuses on only AOC No. 19, as a follow up to the previous investigation recently conducted in August 2011 for all three AOCs in Group 6.

The purpose of this site investigation is to evaluate the presence of chromium recently found in soils at AOC No. 19. This is consistent with the overall goal to determine and evaluate the presence, nature, and extent of releases of contaminants in accordance with 20.4.1.500 New

Mexico Administrative Code (NMAC) incorporating 40 Code of Federal Regulations (CFR) Section 264.101. The investigation activities will be conducted in accordance with Section IV of the Order.

## **Scope of Services**

### **Anticipated Activities**

Pursuant to Section IV of the Order, a scope of services has been developed to evaluate the presence of chromium in soils at AOC No. 19. To accomplish this objective, surface soil samples will be collected at AOC No. 19 (Seep North of MW-45). Hand augured soil borings will be installed at the specified locations along the arroyos (drainage pathways) and samples will be collected from 0-6" below grade surface.

Three soil samples are proposed at locations in the drainage pathways where the highest chromium concentrations were found in soil samples collected in August 2011 [AOC 19-13 (0-0.5'), AOC 19-14 (0-0.5'), and AOC 19-18 (0-0.5')] and at three additional locations within AOC No. 19 where contaminants were not detected at high concentration during the recent investigation.

Soil samples will be collected at AOC No. 19 from the following locations:

- One surface soil sample will be collected near Seep No. 4 at the center of the natural basin where the seep daylighted and previous sample AOC 19-13 was collected;
- One surface soil sample will be collected near Seep No. 9 from along the margin of the drainage pathway where previous sample AOC 19-14 was collected;
- One surface soil sample will be collected from the center of the major drainage pathway between Seep No. 5 and Seep No. 9 where previous sample AOC 19-18 was collected; and
- Three surface soil samples will be collected from locations east of the major drainage pathway that extends from Seep No. 5 to Seep No. 9.

The soil samples will be analyzed for chromium using EPA methods 6010B and 3060A/7196A. EPA method 3060A/7196A is specifically designed to determine the presence and concentration of chromium VI in soils/sediments. EPA method 6010B reports the total chromium concentration (e.g., chromium III plus chromium VI).



## Background Information Research

Documents containing the results of previous investigations<sup>1</sup> and routine groundwater monitoring data<sup>2</sup> from monitoring wells and the seeps were reviewed to facilitate development of this work plan. The previously collected data provide detailed information on the overall subsurface conditions, including hydrogeology and contaminant distribution within groundwater on a site-wide basis. The data collected under this scope of services will supplement the existing soil and groundwater information and provide specific information regarding contaminant occurrence and distribution within soils near the seeps.

## Collection and Management of Investigation Derived Waste

It is not anticipated that excess soil cuttings will be generated since only surface samples (0-0.5') will be collected as part of this Work Plan. All decontamination water will be disposed in the refinery wastewater treatment system upstream of the API Separator.

## Surveys

The horizontal coordinates and elevation of each sampling location and the locations of all other pertinent structures will be determined by a registered New Mexico professional land surveyor in accordance with the State Plane Coordinate System (NMSA 1978 47-1-49-56 (Repl. Pamph. 1993)). The surveys will be conducted in accordance with Sections 500.1 through 500.12 of the Regulations and Rules of the Board of Registration for Professional Engineers and Surveyors Minimum Standards for Surveying in New Mexico. Horizontal positions will be measured to the nearest 0.1-ft and vertical elevations will be measured to the nearest 0.01-ft.

## **Investigation Methods**

### Soil Boring Installation

Shallow soil borings will be completed using a hand auger and/or shovel at appropriate locations within and into the banks of the drainage pathways. This sampling method is ideally suited due to the inaccessibility of a drill rig and the limited anticipated depth to bedrock in the vicinity of the seeps. The sampling equipment will be properly decontaminated before beginning each boring.

---

<sup>1</sup> RPS, 2011, Investigation Report Group 6

<sup>2</sup> Western Refining, 2012, Groundwater Remediation and Monitoring Annual Report

The NMED will be notified as early as practicable if conditions arise or are encountered that do not allow the advancement of borings at planned sampling locations. The sampling will be accomplished under the direction of a qualified engineer or geologist who will maintain a detailed log of the materials and conditions encountered in each boring. Both sample information and visual observations of the samples will be recorded on the boring log. Known site features and/or site survey grid markers will be used as references to locate each boring prior to surveying the location as described above. The boring locations will be measured to the nearest foot, and locations will be recorded on a scaled site map upon completion of each boring.

### Soil Sampling

The individual sample locations have been selected in an attempt to determine the valence state of chromium, which was previously detected in soil samples collected from AOC No. 19. The surface samples may be collected using decontaminated, hand-held stainless steel sampling device, hand auger, or a pre-cleaned disposable sampling device. A portion of the sample will be placed in pre-cleaned, laboratory-prepared sample containers for laboratory chemical analysis. The remaining portions of the sample will be used for logging and field screening as discussed below. Sample handling and chain-of-custody procedures will be in accordance with the procedures presented below.

Discrete soil samples will be collected from 0-6" (all soil sample locations). Three soil samples are proposed at locations in the drainage pathways where the highest chromium concentrations were found in soil samples collected in August 2011 [AOC 19-13 (0-0.5'), AOC 19-14 (0-0.5'), and AOC 19-18 (0-0.5')] and at three additional locations within AOC No. 19 where contaminants were not detected at high concentration during the recent Group 6 investigation. The proposed sample locations are shown on enclosed Figure 2.

Quality Assurance/Quality Control (QA/QC) samples will be collected to monitor the validity of the soil sample collection procedures as follows:

- One field duplicate will be collected and analyzed for total chromium.

### Soil Sample Field Screening and Logging

Samples will be screened in the field for evidence of contaminants. Field screening results will be recorded on the exploratory boring logs. The primary screening methods include: (1) visual examination, and (2) olfactory examination.

Visual screening includes examination of soil samples for evidence of staining caused by petroleum-related compounds. Field screening results may be site- and boring-specific and the results may vary with instrument type, the media screened, weather conditions, moisture content, soil type, and type of contaminant. Conditions capable of influencing the results of field screening will be recorded on the field logs.

The physical characteristics of the samples (such as mineralogy, ASTM soil classification, moisture content, texture, color, presence of stains or odors, and/or field screening results), depth where each sample was obtained, method of sample collection, and other observations will be recorded in the field log by a qualified geologist or engineer. Detailed logs of each soil boring will be completed in the field by a qualified engineer or geologist. Additional information, such as any unusual or noticeable conditions encountered during sampling, will be recorded on the logs.

### Sample Handling

At a minimum, the following procedures will be used at all times when collecting samples during investigation, corrective action, and monitoring activities:

1. Neoprene, nitrile, or other protective gloves will be worn when collecting samples. New disposable gloves will be used to collect each sample;
2. All samples collected of each medium for chemical analysis will be transferred into clean sample containers supplied by the project analytical laboratory with the exception of soil, rock, and sediment samples obtained in Encore® samplers. Sample container volumes and preservation methods will be in accordance with the most recent standard EPA and industry accepted practices for use by accredited analytical laboratories. Sufficient sample volume will be obtained for the laboratory to complete the method-specific QC analyses on a laboratory-batch basis; and
3. Sample labels and documentation will be completed for each sample following procedures discussed below. Immediately after the samples are collected, they will be stored in a cooler with ice or other appropriate storage method until they are delivered to the analytical laboratory. Standard chain-of-custody procedures, as described below, will be followed for all samples collected. All samples will be submitted to the laboratory soon enough to allow the laboratory to conduct the analyses within the method holding times. At a minimum, all samples will be submitted to the laboratory within 48 hours after their collection.

Chain-of-custody and shipment procedures will include the following:

1. Chain-of-custody forms will be completed at the end of each sampling day, prior to the transfer of samples off site.
2. Individual sample containers will be packed to prevent breakage and transported in a sealed cooler with ice or other suitable coolant or other EPA or industry-wide accepted

method. The drainage hole at the bottom of the cooler will be sealed and secured in case of sample container leakage.

3. Each cooler or other container will be delivered directly to the analytical laboratory.
4. Glass bottles will be separated in the shipping container by cushioning material to prevent breakage.
5. Plastic containers will be protected from possible puncture during shipping using cushioning material.
6. The chain-of-custody form and sample request form will be shipped inside the sealed storage container to be delivered to the laboratory.
7. Chain-of-custody seals will be used to seal the sample-shipping container in conformance with EPA protocol.
8. Signed and dated chain-of-custody seals will be applied to each cooler prior to transport of samples from the site.
9. Upon receipt of the samples at the laboratory, the custody seals will be broken, the chain-of-custody form will be signed as received by the laboratory, and the conditions of the samples will be recorded on the form. The original chain-of-custody form will remain with the laboratory and copies will be returned to the relinquishing party.
10. Copies of all chain-of-custody forms generated as part of sampling activities will be maintained on-site.

### Decontamination Procedures

The objective of the decontamination procedures is to minimize the potential for cross-contamination. Sampling or measurement equipment, including but not limited to, stainless steel sampling tools will be decontaminated in accordance with the following procedures or other methods approved by the Department before each sampling attempt or measurement:

1. Brush equipment with a wire or other suitable brush, if necessary or practicable, to remove large particulate matter;
2. Rinse with potable tap water;
3. Wash with nonphosphate detergent or other detergent approved by the Department (examples include Fantastik™, Liqui-Nox®);
4. Rinse with potable tap water; and
5. Double rinse with deionized water.

All decontamination solutions will be collected and stored temporarily as described above in the discussion on the collection and management of investigation derived waste. Decontamination procedures and the cleaning agents used will be documented in the daily field log.

### Documentation of Field Activities

Daily field activities, including observations and field procedures, will be recorded in a field log book. The original field forms will be maintained at the facility. Copies of the completed forms will be maintained in a bound and sequentially numbered field file for reference during field activities. Indelible ink will be used to record all field activities. Photographic documentation of field activities will be performed, as appropriate. The daily record of field activities will include the following:

1. Site or unit designation;
2. Date;
3. Time of arrival and departure;
4. Field investigation team members including subcontractors and visitors;
5. Weather conditions;
6. Daily activities and times conducted;
7. Observations;
8. Record of samples collected with sample designations and locations specified;
9. Photographic log, as appropriate;
10. Field monitoring data, including health and safety monitoring;
11. Equipment used and calibration records, if appropriate;
12. List of additional data sheets and maps completed;
13. An inventory of the waste generated and the method of storage or disposal; and
14. Signature of personnel completing the field record.

### Chemical Analyses

All samples collected for laboratory analysis will be submitted to an accredited laboratory. The laboratory will use the most recent standard EPA and industry-accepted analytical methods for target analytes as the testing methods for each medium sampled. Chemical analyses will be performed in accordance with the most recent EPA standard analytical methodologies and extraction methods.

The soil samples will be analyzed by the following methods:

- SW-846 Method 6010B for total chromium; and
- SW-846 Method 3060A/7196A for chromium VI.

## **Schedule**

This Investigation Work Plan will be implemented concurrently with the Group 9 Investigation Work Plan pending NMED approval. The estimated timeframes for each of the planned activities is as shown below:

- Field work (inclusive of all soil sampling) – one week;
- Laboratory analyses – four weeks;
- Data reduction and validation – two weeks; and
- Data gap analysis – two weeks.

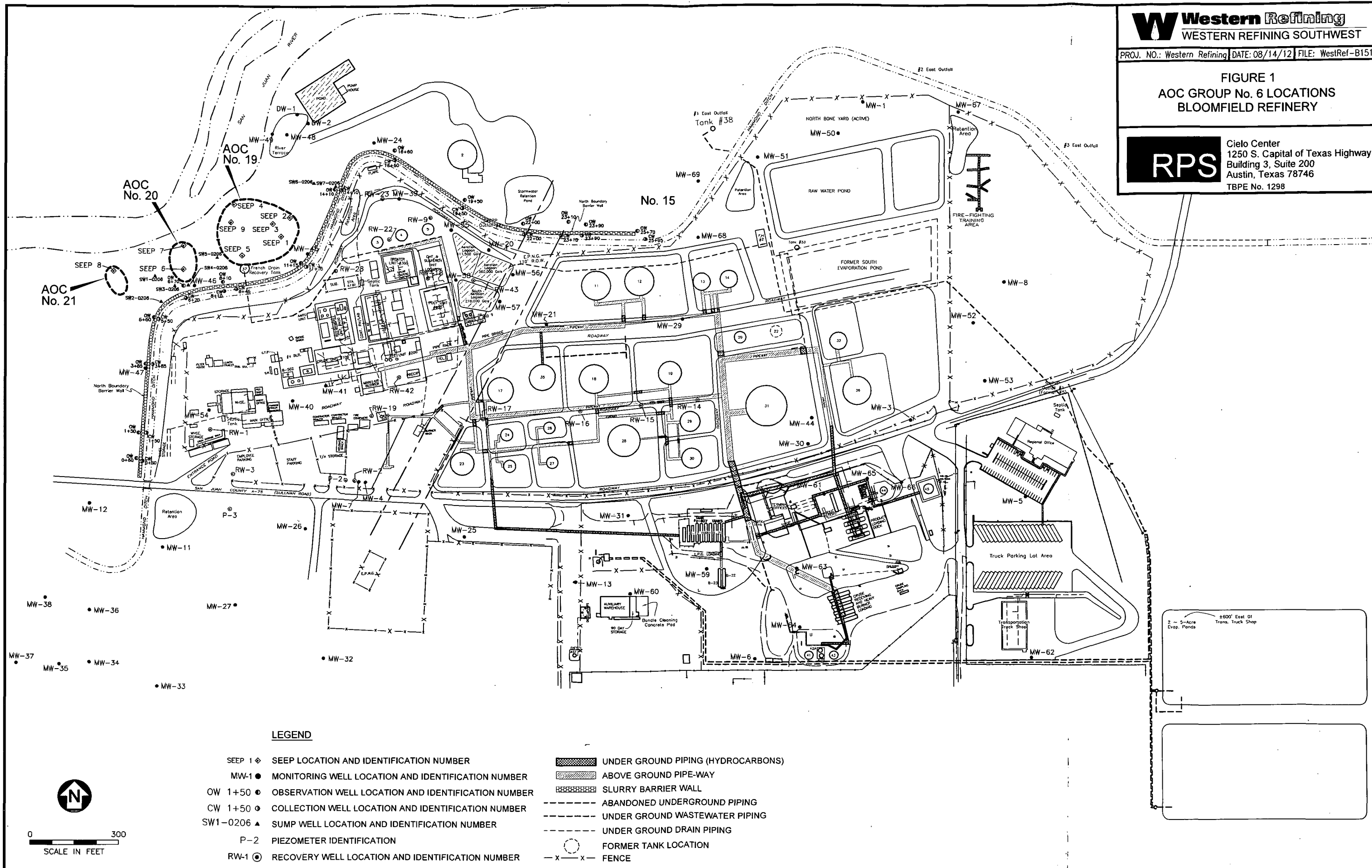
Completion of the data gap analysis will complete all activities conducted under this Investigation Work Plan.

## FIGURES

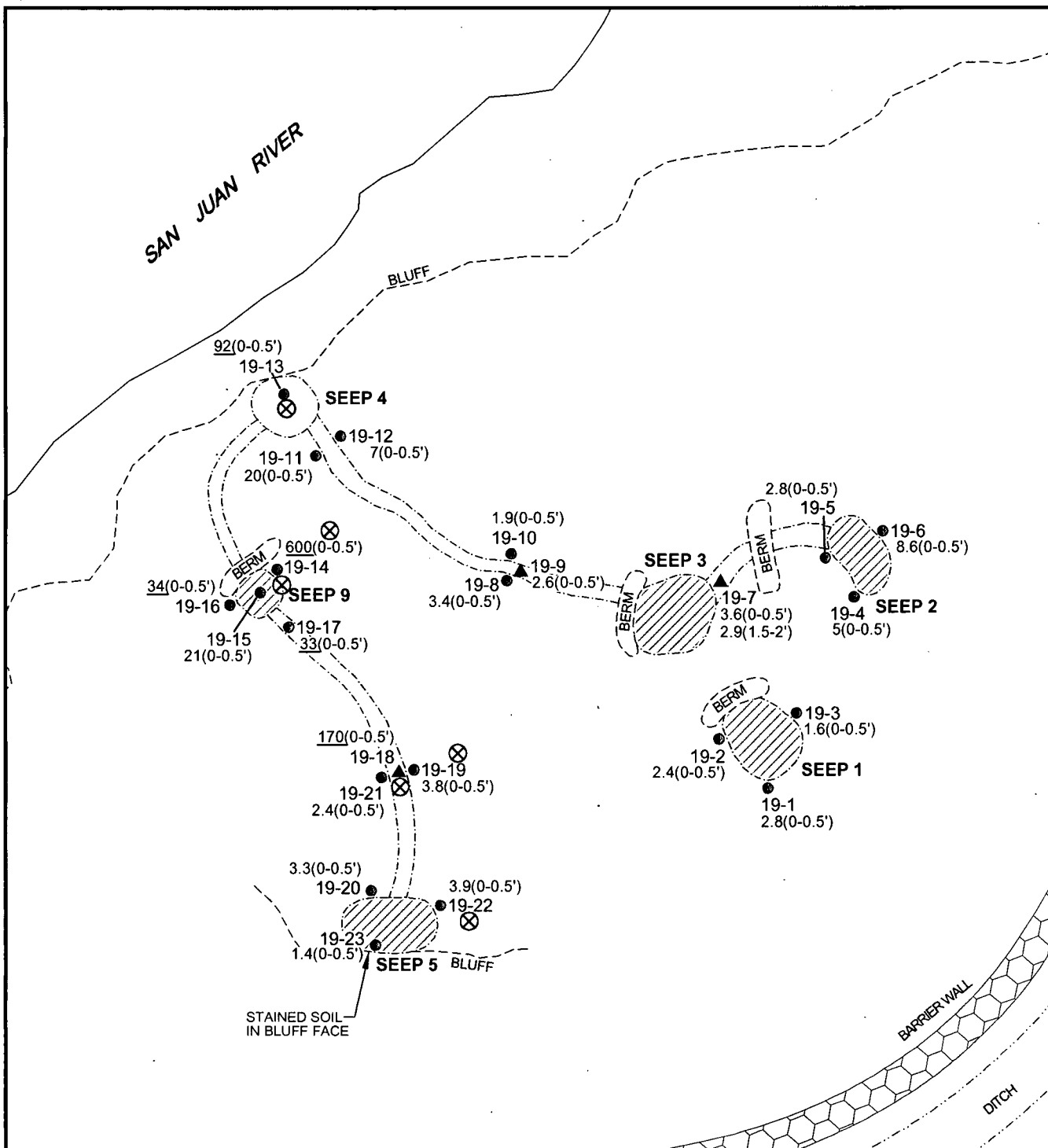
FIGURE 1  
AOC GROUP No. 6 LOCATIONS  
BLOOMFIELD REFINERY

**RPS**

Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298







#### LEGEND

- ⊗ PROPOSED NEW SOIL SAMPLE LOCATION
- SEEP 1** SEEP IDENTIFICATION NUMBER
- 19-1 ● SURFACE SOIL SAMPLE LOCATION AND IDENTIFICATION NUMBER
- 19-7 ▲ SOIL BORING LOCATION AND IDENTIFICATION NUMBER
- ▨ LINED CATCHMENT BASIN
- DRAINAGE PATHWAY
- 2.4(0-0.5') CHROMIUM CONCENTRATION, mg/kg (SAMPLE DEPTH - FT)
- 23.7 CONCENTRATION EXCEEDS DAF SCREENING LEVEL



0 50  
SCALE IN FEET

**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 08/14/12 FILE: WestRef-A137

#### FIGURE 2 AOC No. 19 CHROMIUM SOIL MAP BLOOMFIELD REFINERY

**RPS**

Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298

July 26, 2012

RECEIVED OCD

2012 JUL 30 P 12: 55

John E. Kieling, Chief  
New Mexico Environment Department  
Hazardous Waste Bureau  
2905 Rodeo Park Drive East, Bldg 1  
Santa Fe, NM 87505

**UPS Tracking #: 1Z 881 839 24 5595 6248 (Delivery to NMED)**  
**Certified Mail #: 7009 2250 0002 3833 5315 (Delivery to OCD)**

**RE:** Response to NMED's Approval  
Investigation Report Group 6 (AOC No. 19 Seep North of MW-45, AOC No. 20 Seep North of MW-46, and AOC No. 21 Seep North of MW-47)  
Western Refining Southwest, Inc., Bloomfield Refinery  
EPA ID# NMD089416416  
HWB-WRB-11-008

Dear Mr. Kieling:

Pursuant to your approval letter dated July 5, 2012, Western Refining Southwest, Inc., Bloomfield Refinery ("Western") has prepared the following responses to address your comments. Please find the requested revisions enclosed and note that the requested investigation work plan for Group 6 will be submitted on or before October 12, 2012, as requested.

**NMED Comment No. 1:**

*Section 7 (Conclusions and Recommendations), pages 32 through 34, summarizes the results of the investigation and provides conclusions and recommendations for the work completed at Group 6. However, Western did not discuss or provide any recommendations to address the DRO result of 21,000 mg/kg from soil sample location AOC 19-23, which is significantly higher than the residential soil screening level (SSLs). NMED understands that the conditions at the site are not conducive to allow removal of the material from the bluff surface; however, Western must provide an explanation about the conditions at soil sample location AOC 19-23 to include a description of the location and an explanation why removal is not warranted or practicable. Western must include this discussion at the end of Section 7.2 (Recommendations) and provide a replacement page for the Report.*

**Western Response No. 1:** Pursuant to our earlier discussions, please find enclosed replacement pages for Section 7. The text has been revised to clarify the actual site conditions in the area of sample location AOC 19-23. A redline-strikeout version is also included to facilitate your review.

**NMED Comment No. 2:**

*Section 7.2 (Recommendations), page 34, states that "[a]dditional soil samples should be collected at the locations with the highest concentrations of chromium at AOC 19 [e.g., AOC 19-13 (0-0.5'), AOC 19-14 (0-0.5'), and AOC 19-18 (0-0.5')]. The samples should be analyzed for chromium VI to determine if the chromium detected in soils at AOC 19 is present as chromium III or chromium VI." Western must collect at least three additional soil samples from*

*soils within AOC 19 at locations where contaminants are not detected at high concentrations to determine if chromium is typically present as chromium III and not present only as a result of local reducing conditions in areas containing petroleum contamination. Western must submit a letter work plan to propose the additional investigation activities to collect these soil samples.*

**Western Response No. 2:** A work plan for collection of additional soil samples for chromium analysis will be submitted, as directed by NMED, on or before October 12, 2012.

**NMED Comment No. 3:**

*Figures 12 (AOC No. 19 Arsenic Soil Map) through 19 (AOC No. 21 Diesel Range Organics Soil Map) depict surface soil and soil boring locations with their corresponding analytical results. In addition, the results greater than the SSL are underlined. However, Western did not specify which SSL is represented in the figures. In future submittals, specify which SSLs are referenced in figures.*

**Western Response No. 3:** In future submittals, the SSL value will be explained in the legend of any associated figures.

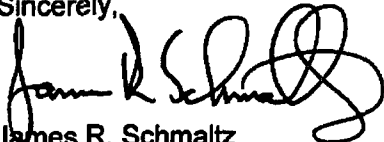
**NMED Comment No. 4:**

*Appendix C provides site photographs of the areas of concern (AOCs) and seep locations. Some of the site photos depict sample locations, but they are not identified in the captions. In future submittals, identify the sample locations in the captions below the site photos if the sample locations are visible in the photo.*

**Western Response No. 4:** In future submittals, any sample locations that appear in site photos will be identified in a caption below the photo.

If you have questions regarding the above responses or the enclosure, please feel free to contact me at (505) 632-4171 at your convenience.

Sincerely,



James R. Schmaltz  
Health, Safety, Environmental and Regulatory Director  
Western Refining Southwest, Inc., Bloomfield Refinery

Enclosure

cc: Dave Cobrain – NMED HWB  
Leona Tsinnajinnie – NMED HWB  
Carl Chavez - NMOCD  
Allen Hains – Western Refining El Paso

## Section 7

# Conclusions and Recommendations

This section summarizes and provides an evaluation of the potential impacts as shown in field screening data and analytical data. An investigation of soils was conducted at Group 6 (AOCs 19, 20, and 21) to assess and evaluate the presence, nature, extent, fate, and transport of contaminants. To accomplish this objective, soil samples were collected and analyzed for potential site-related constituents.

### 7.1 Conclusions

#### AOC No. 19 – Seep North of MW-45

Three metals (arsenic, cobalt, and chromium) and diesel range organics were detected in soil samples at concentrations exceeding their respective screening levels. Arsenic was detected in nine samples at concentrations above the DAF screening level, with eight of these samples having concentrations above the residential screening level. Cobalt was detected in generally low concentrations with a maximum detected value of 11 mg/kg. As shown in Table 5, the samples with concentrations of cobalt above the screening levels do not have concentrations of organic constituents above screening levels. Chromium was detected in five samples at concentrations above the DAF screening level with only one being above the residential screening level. Arsenic, chromium, and cobalt are naturally occurring constituents in many soils and it is possible that these constituents are naturally occurring in soils at the Bloomfield Refinery. The analyses for chromium were for total chromium, which are compared to the chromium VI screening levels. It may be overly conservative to compare to the total chromium results to the chromium VI screening levels, but additional analyses will be required to speciate the chromium.

The only analytical results for organic constituents (i.e., diesel range organics) above only the DAF screening level occur at three locations with a maximum value of 770 mg/kg. All three of these detections of diesel range organics are below the residential screening level. A fourth sample (AOC 19-23), which was collected from a cliff face, had a concentration of diesel range organics above the DAF and residential screening levels. This single detection of diesel range organics above the residential screening level does not warrant a response action because there is very limited potential for exposure to the small area represented by this sample.

The AOC 19-23 sample was collected from the location on a cliff face that had the greatest visible staining from apparent hydrocarbons. The entire cliff face at this location is only approximately five feet high and seven feet long, and there is very limited access to this area. The seep that caused the presence of hydrocarbons is no longer active with the installation of the slurry wall, therefore the hydrocarbons present at the surface should naturally degrade over time. The actual volume of impacted soils present near the land surface is also very small, as it appears the previous seep mainly caused only surface staining of soils along the small cliff face.

#### AOC No. 20 – Seep North of MW-46

Only cobalt and diesel range organics were detected at concentrations above the applicable screening levels. Cobalt was detected in only one soil sample at a concentration (8.9 mg/kg) above the DAF screening level (5.51 mg/kg) but did not exceed the residential soil screening level of 23 mg/kg. As noted above, cobalt may be a naturally occurring constituent. Diesel range organics were detected in only one sample at a concentration (490 mg/kg) above the DAF soil screening level of 200 mg/kg. None of the diesel range organic analytical results exceeded the residential direct contact screening level of 1,830 mg/kg.

#### AOC No. 21 – Seep North of MW-47

Similarly to AOC No. 20, only cobalt and diesel range organics were detected at concentrations above the applicable screening levels. Cobalt was detected in eight soil samples at concentrations above the DAF screening level (5.51 mg/kg) but none exceed the residential soil screening level of 23 mg/kg. Diesel range organics were detected in one sample at a concentration (710 mg/kg) above the DAF soil screening level of 200 mg/kg. None of the diesel range organic analytical results exceed the residential direct contact screening level of 1,830 mg/kg.

#### Cumulative Risk Evaluation

A cumulative risk evaluation for soils is presented in Table 6. This was conducted by taking the maximum reported concentration of each detected constituent from all three AOCs and dividing by the residential screening level and non-residential screening levels as shown in the equations below. These calculations are separated for carcinogenic and non-carcinogenic constituents. The cumulative carcinogenic risk is  $6.703 \times 10^{-5}$  assuming residential land use

and  $1.36 \times 10^{-5}$  for non-residential land use. The hazard index for residential land use is 3.557 and for non-residential land use is 1.652.

$$\text{Site Risk} = \left( \frac{\text{conc}_x}{\text{SSL}_x} + \frac{\text{conc}_y}{\text{SSL}_y} + \frac{\text{conc}_z}{\text{SSL}_z} + \dots + \frac{\text{conc}_i}{\text{SSL}_i} \right) \times 10^{-5}$$

$$\text{Site Hazard Index (HI)} = \left( \frac{\text{conc}_x}{\text{SSL}_x} + \frac{\text{conc}_y}{\text{SSL}_y} + \frac{\text{conc}_z}{\text{SSL}_z} + \dots + \frac{\text{conc}_i}{\text{SSL}_i} \right) \times 1$$

## 7.2 Recommendations

Western plans to conduct an investigation of background locations and will attempt to establish site-specific background values for various naturally occurring constituents (e.g., arsenic and cobalt in soils). After background values have been established, site data collected for AOCs 19, 20 and 21 should be reassessed to determine if some analytical results are actually representative of background conditions and not site-related impacts.

Additional soil samples should be collected at the locations with the highest concentrations of chromium at AOC 19 [e.g., AOC19-13 (0-0.5'), AOC 19-14 (0-0.5'), and AOC 19-18 (0-0.5')]. The samples should be analyzed for chromium VI to determine if the chromium detected in soils at AOC 19 is present as chromium III or chromium VI. If the chromium is determined to be chromium III instead of chromium VI, then the screening levels in Table 5 will need to be revised along with any associated figures.

After site-specific background values have been established and the location-specific speciation of chromium has been determined, then preparation of a risk assessment for the Group 6 AOCs may be appropriate. The site concentrations have been compared to residential screening levels in this investigation report; however, the physical nature of the seep areas may limit future use of this area. An evaluation of potential exposures in a site-specific risk assessment could offer a more appropriate evaluation of actual risks based on more likely future use of the area at AOCs 19, 20, and 21.

## **Section 7**

# **Conclusions and Recommendations**

This section summarizes and provides an evaluation of the potential impacts as shown in field screening data and analytical data. An investigation of soils was conducted at Group 6 (AOCs 19, 20, and 21) to assess and evaluate the presence, nature, extent, fate, and transport of contaminants. To accomplish this objective, soil samples were collected and analyzed for potential site-related constituents.

### **7.1 Conclusions**

#### AOC No. 19 – Seep North of MW-45

Three metals (arsenic, cobalt, and chromium) and diesel range organics were detected in soil samples at concentrations exceeding their respective screening levels. Arsenic was detected in nine samples at concentrations above the DAF screening level, with eight of these samples having concentrations above the residential screening level. Cobalt was detected in generally low concentrations with a maximum detected value of 11 mg/kg. As shown in Table 5, the samples with concentrations of cobalt above the screening levels do not have concentrations of organic constituents above screening levels. Chromium was detected in five samples at concentrations above the DAF screening level with only one being above the residential screening level. Arsenic, chromium, and cobalt are naturally occurring constituents in many soils and it is possible that these constituents are naturally occurring in soils at the Bloomfield Refinery. The analyses for chromium were for total chromium, which are compared to the chromium VI screening levels. It may be overly conservative to compare to the total chromium results to the chromium VI screening levels, but additional analyses will be required to speciate the chromium.

The analytical results for organic constituents (i.e., diesel range organics) above only the DAF screening level occur at three locations with a maximum value of 770 mg/kg. All three of these detections of diesel range organics are below the residential screening level. A fourth sample (AOC 19-23), which was collected from a cliff face, had a concentration of diesel range organics above the DAF and residential screening levels. This single detection of diesel range organics above the residential screening level does not warrant a response action because there is very limited potential for exposure to the small area represented by this sample.

The AOC 19-23 sample was collected from the location on a cliff face that had the greatest visible staining from apparent hydrocarbons. The entire cliff face at this location is only approximately five feet high and seven feet long, and there is very limited access to this area. The seep that caused the presence of hydrocarbons is no longer active with the installation of the slurry wall, therefore the hydrocarbons present at the surface should naturally degrade over time. The actual volume of impacted soils present near the land surface is also very small, as it appears the previous seep mainly caused only surface staining of soils along the small cliff face.

#### AOC No. 20 – Seep North of MW-46

Only cobalt and diesel range organics were detected at concentrations above the applicable screening levels. Cobalt was detected in only one soil sample at a concentration (8.9 mg/kg) above the DAF screening level (5.51 mg/kg) but did not exceed the residential soil screening level of 23 mg/kg. As noted above, cobalt may be a naturally occurring constituent. Diesel range organics were detected in only one sample at a concentration (490 mg/kg) above the DAF soil screening level of 200 mg/kg. None of the diesel range organic analytical results exceeded the residential direct contact screening level of 1,830 mg/kg.

#### AOC No. 21 – Seep North of MW-47

Similarly to AOC No. 20, only cobalt and diesel range organics were detected at concentrations above the applicable screening levels. Cobalt was detected in eight soil samples at concentrations above the DAF screening level (5.51 mg/kg) but none exceed the residential soil screening level of 23 mg/kg. Diesel range organics were detected in one sample at a concentration (710 mg/kg) above the DAF soil screening level of 200 mg/kg. None of the diesel range organic analytical results exceed the residential direct contact screening level of 1,830 mg/kg.

#### Cumulative Risk Evaluation

A cumulative risk evaluation for soils is presented in Table 6. This was conducted by taking the maximum reported concentration of each detected constituent from all three AOCs and dividing by the residential screening level and non-residential screening levels as shown in the equations below. These calculations are separated for carcinogenic and non-carcinogenic constituents. The cumulative carcinogenic risk is  $6.703 \times 10^{-5}$  assuming residential land use



and  $1.36 \times 10^{-5}$  for non-residential land use. The hazard index for residential land use is 3.557 and for non-residential land use is 1.652.

$$\text{Site Risk} = \left( \frac{\text{conc}_x}{\text{SSL}_x} + \frac{\text{conc}_y}{\text{SSL}_y} + \frac{\text{conc}_z}{\text{SSL}_z} + \dots + \frac{\text{conc}_i}{\text{SSL}_i} \right) \times 10^{-5}$$

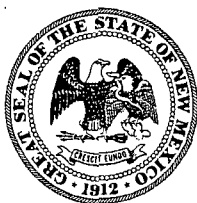
$$\text{Site Hazard Index (HI)} = \left( \frac{\text{conc}_x}{\text{SSL}_x} + \frac{\text{conc}_y}{\text{SSL}_y} + \frac{\text{conc}_z}{\text{SSL}_z} + \dots + \frac{\text{conc}_i}{\text{SSL}_i} \right) \times 1$$

## 7.2 Recommendations

Western plans to conduct an investigation of background locations and will attempt to establish site-specific background values for various naturally occurring constituents (e.g., arsenic and cobalt in soils). After background values have been established, site data collected for AOCs 19, 20 and 21 should be reassessed to determine if some analytical results are actually representative of background conditions and not site-related impacts.

Additional soil samples should be collected at the locations with the highest concentrations of chromium at AOC 19 [e.g., AOC19-13 (0-0.5'), AOC 19-14 (0-0.5'), and AOC 19-18 (0-0.5')]. The samples should be analyzed for chromium VI to determine if the chromium detected in soils at AOC 19 is present as chromium III or chromium VI. If the chromium is determined to be chromium III instead of chromium VI, then the screening levels in Table 5 will need to be revised along with any associated figures.

After site-specific background values have been established and the location-specific speciation of chromium has been determined, then preparation of a risk assessment for the Group 6 AOCs may be appropriate. The site concentrations have been compared to residential screening levels in this investigation report; however, the physical nature of the seep areas may limit future use of this area. An evaluation of potential exposures in a site-specific risk assessment could offer a more appropriate evaluation of actual risks based on more likely future use of the area at AOCs 19, 20, and 21.



SUSANA MARTINEZ  
Governor

JOHN A. SANCHEZ  
Lieutenant Governor

NEW MEXICO  
ENVIRONMENT DEPARTMENT

*Hazardous Waste Bureau*

2905 Rodeo Park Drive East, Building 1  
Santa Fe, New Mexico 87505-6303  
Phone (505) 476-6000 Fax (505) 476-6030  
[www.nmenv.state.nm.us](http://www.nmenv.state.nm.us)



DAVE MARTIN  
Secretary

BUTCH TONGATE  
Deputy Secretary

JAMES H. DAVIS, Ph.D.  
Director  
Resource Protection Division

CERTIFIED MAIL - RETURN RECEIPT REQUESTED

July 5, 2012

Mr. Randy Schmaltz  
Environmental Manager  
Western Refining, Bloomfield Refinery  
P.O. Box 159  
Bloomfield, New Mexico 87413

**RE: APPROVAL  
INVESTIGATION REPORT GROUP 6 (AOC NO. 19 SEEP  
NORTH OF MW-45, AOC NO. 20 SEEP NORTH OF MW-46, AND  
AOC NO. 21 SEEP NORTH OF MW-47)  
WESTERN REFINING SOUTHWEST, INC., BLOOMFIELD REFINERY  
EPA ID# NMD089416416  
HWB-WRB-11-008**

Dear Mr. Schmaltz:

The New Mexico Environment Department (NMED) has completed its review of Western Refining Southwest, Inc., Bloomfield Refinery's (Western) *Investigation Report Group 6 (AOC No. 19 Seep North of MW-45, AOC No. 20 Seep North of MW-46, and AOC No. 21 Seep North of MW-47)* (Report), dated November 2011. NMED hereby approves this Report with the following comments.

**Comment 1**

Section 7 (Conclusions and Recommendations), pages 32 through 34, summarizes the results of the investigation and provides conclusions and recommendations for the work completed at Group 6. However, Western did not discuss or provide any recommendations to address the

R. Schmaltz

July 5, 2012

Page 2 of 3

DRO result of 21,000 mg/kg from soil sample location AOC 19-23, which is significantly higher than the residential soil screening level (SSLs). NMED understands that the conditions at the site are not conducive to allow removal of the material from the bluff surface; however, Western must provide an explanation about the conditions at soil sample location AOC 19-23 to include a description of the location and an explanation why removal is not warranted or practicable. Western must include this discussion at the end of Section 7.2 (Recommendations) and provide a replacement page for the Report.

### **Comment 2**

Section 7.2 (Recommendations), page 34, states that “[a]dditional soil samples should be collected at the locations with the highest concentrations of chromium at AOC 19 [e.g., AOC 19-13 (0-0.5’), AOC 19-14 (0-0.5’), and AOC 19-18 (0-0.5’)]. The samples should be analyzed for chromium VI to determine if the chromium detected in soils at AOC 19 is present as chromium III or chromium VI.” Western must collect at least three additional soil samples from soils within AOC 19 at locations where contaminants are not detected at high concentrations to determine if chromium is typically present as chromium III and not present only as a result of local reducing conditions in areas containing petroleum contamination. Western must submit a letter work plan to propose the additional investigation activities to collect these soil samples.

### **Comment 3**

Figures 12 (AOC No. 19 Arsenic Soil Map) through 19 (AOC No. 21 Diesel Range Organics Soil Map) depict surface soil and soil boring locations with their corresponding analytical results. In addition, the results greater than the SSL are underlined. However, Western did not specify which SSL is represented in the figures. In future submittals, specify which SSLs are referenced in figures.

### **Comment 4**

Appendix C provides site photographs of the areas of concern (AOCs) and seep locations. Some of the site photos depict sample locations, but they are not identified in the captions. In future submittals, identify the sample locations in the captions below the site photos if the sample locations are visible in the photo.

R. Schmaltz  
July 5, 2012  
Page 3 of 3

Western must adhere to all requirements in this Approval in future submittals. The replacement page must be submitted to NMED by **July 27, 2012**. The work plan referenced in Comment 2 must be submitted to NMED no later than **October 12, 2012**.

If you have any questions regarding this letter, please contact Leona Tsinnajinnie of my staff at (505) 476-6057.

Sincerely,

A handwritten signature in dark ink, appearing to read "John E. Kielling" with a stylized flourish at the end.

John E. Kielling  
Chief  
Hazardous Waste Bureau

cc: D. Cobrain, NMED HWB  
C. Chavez, OCD  
A. Hains, Western  
K. Robinson, Western Bloomfield Refinery

File: HWB-WRB-11-008 and Reading 2012

November 11, 2011

John E. Kieling, Acting Bureau Chief  
New Mexico Environment Department  
Hazardous Waste Bureau  
2905 Rodeo Park Drive East, Bldg 1  
Santa Fe, NM 87505

UPS Mail #: 1ZF9F6470196574075 (Addressed to NMED)  
Certified Mail #: 7009 2250 0002 3833 5094 (Addressed to OCD)

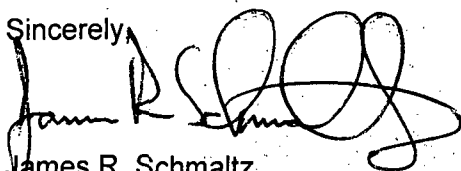
**Re: Group 6 Investigation Report  
Western Refining Southwest, Inc., Bloomfield Refinery  
EPA ID# NMD089416416**

Dear Mr. Kieling:

Western Refining Southwest, Inc., Bloomfield Refinery submits the referenced Investigation Report pursuant to Section IV.B.7 of the July 2007 HWB Order. This Investigation Report summarizes the site environmental investigation activities completed for AOC No. 19 Seep North of MW-45, AOC No. 20 Seep North of MW-46, and AOC No. 21 Seep North of MW-47, which has been designated as Group 6.

If you have any questions or would like to discuss the Investigation Report, please contact me at (505) 632-4171

Sincerely,



James R. Schmaltz  
Health, Safety, Environmental, and Regulatory Director  
Western Refining Southwest, Inc.  
Bloomfield Refinery

cc: Dave Cobrain – NMED HWB  
Leona Tsinnajinnie – NMED HWB  
Carl Chavez - NMOCD  
Allen Hains – Western Refining El Paso  
Scott Crouch – RPS Austin



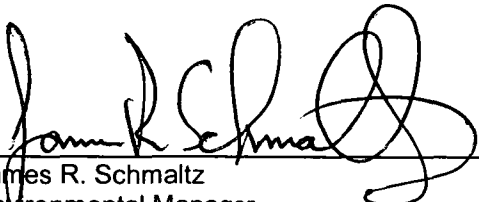
Cielo Center, 1250 South Capital of Texas Highway, Building Three, Suite 200, Austin, Texas 78746, USA  
T +1 512 347 7588 F +1 512 347 8243 W [www.rpsgroup.com](http://www.rpsgroup.com)

## INVESTIGATION REPORT GROUP 6

(AOC No. 19 Seep North of MW-45,  
AOC No. 20 Seep North of MW-46,  
and AOC No. 21 Seep North of MW-47)

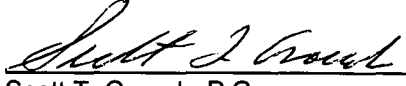
**Bloomfield Refinery**  
**Western Refining Southwest, Inc.**  
**#50 Rd 4990**  
**Bloomfield, New Mexico 87413**

**November 2011**



---

James R. Schmaltz  
Environmental Manager  
Western Refining Southwest, Inc.  
Bloomfield Refinery



---

Scott T. Crouch, P.G.  
Senior Consultant  
RPS  
1250 South Capital of Texas Highway  
Building Three, Suite 200  
Austin, Texas 78746

# Table of Contents

## List of Sections

|  |    |
|--|----|
| List of Acronyms .....   | 1  |
| Executive Summary .....  | 2  |
| Section 1 Introduction .....   | 4  |
| Section 2 Background.....  | 6  |
| Section 3 Scope of Activities .....  | 9  |
| 3.1 Soil Boring Installation and Sample Collection .....                           | 9  |
| 3.2 Background Information Research .....  | 11 |
| 3.3 Collection and Management of Investigation Derived Waste .....                 | 11 |
| 3.4 Surveys .....  | 12 |
| Section 4 Field Investigation Results .....  | 13 |
| 4.1 Surface Conditions .....   | 13 |
| 4.2 Subsurface Conditions.....   | 14 |
| 4.3 Exploratory Drilling Investigations, Soil Sampling and Boring Abandonment..... | 15 |
| 4.3.1 AOC 19 – Seep North of MW-45.....  | 15 |
| 4.3.2 AOC 20 – Seep North of MW-46.....  | 20 |
| 4.3.3 AOC 21 – Seep North of MP-47.....  | 22 |
| 4.4 Groundwater Conditions .....   | 23 |
| 4.5 Surface Water Conditions .....   | 24 |
| Section 5 Regulatory Criteria .....  | 26 |
| Section 6 Site Impacts .....   | 29 |
| 6.1 Soil Sampling Chemical Analytical Results.....                                 | 29 |
| Section 7 Conclusions and Recommendations .....                                    | 32 |
| 7.1 Conclusions .....  | 32 |
| 7.2 Recommendations .....  | 33 |
| Section 8 References .....   | 35 |

## List of Tables

|         |  |
|---------|--|
| Table 1 | Historical Seep Discharge Analyses                       |
| Table 2 | Group 6 Soil Boring Samples – Vapor Screening Results    |
| Table 3 | Residential Soil Screening Levels                        |
| Table 4 | Non-Residential Soil Screening Levels                    |
| Table 5 | Group 6 Soil Analytical Results Summary                  |
| Table 6 | Group 6 Soil Cumulative Risk and Hazard Index Evaluation |

## **Table of Contents (*continued*)**

### **List of Figures**

|           |  |
|-----------|--|
| Figure 1  | Site Location Map  |
| Figure 2  | Group No. 6 AOC Locations                                |
| Figure 3  | AOC Nos. 19, 20 and 21 Seep Location Map                 |
| Figure 4  | AOC No. 19 Sample Location Map                           |
| Figure 5  | AOC No. 20 Sample Location Map                           |
| Figure 6  | AOC No. 21 Sample Location Map                           |
| Figure 7  | Cross Section A-A' West to East                          |
| Figure 8  | Cross Section B-B' North to South                        |
| Figure 9  | Potentiometric Surface Map August 2009                   |
| Figure 10 | Separate Phase Hydrocarbon Thickness Map August 2008     |
| Figure 11 | Dissolved-Phase Groundwater Data August and October 2008 |
| Figure 12 | AOC No. 19 Arsenic Soil Map                              |
| Figure 13 | AOC No. 19 Chromium Soil Map                             |
| Figure 14 | AOC No. 19 Cobalt Soil Map                               |
| Figure 15 | AOC No. 19 Diesel Range Organics Soil Map                |
| Figure 16 | AOC No. 20 Cobalt Soil Map                               |
| Figure 17 | AOC No. 20 Diesel Range Organics Soil Map                |
| Figure 18 | AOC No. 21 Cobalt Soil Map                               |
| Figure 19 | AOC No. 21 Diesel Range Organics Map                     |

### **List of Appendices**

|            |  |
|------------|--|
| Appendix A | Field Methods  |
| Appendix B | Survey Data  |
| Appendix C | Photographs  |
| Appendix D | Boring Logs  |
| Appendix E | Site-Specific Dilution/Attenuation Factor Calculations |
| Appendix F | TPH Screening Level Calculations                       |
| Appendix G | Analytical Data Reports                                |
| Appendix H | Quality Assurance/Quality Control Review               |



# List of Acronyms

American Petroleum Institute (API)  
areas of concern (AOCs)  
benzene, toluene, ethylbenzene, and xylene (BTEX)  
below ground level (bgl)  
Code of Federal Regulations (CFR)  
dilution/attenuation factor (DAF)  
Environmental Protection Agency (EPA)  
hazard index (HI)  
hollow-stem auger (HSA)  
investigation derived waste (IDW)  
liquefied petroleum gas (LPG)  
Massachusetts Department of Environmental Protection (MADEP)  
maximum contaminant level (MCL)  
monitoring well (MW)  
New Mexico Administrative Code (NMAC)  
New Mexico Environment Department (NMED)  
Resource Conservation and Recovery Act (RCRA)  
photoionization detector (PID)  
polyvinyl chloride (PVC)  
separate phase hydrocarbon (SPH)  
Solid Waste Management Units (SWMUs)  
total petroleum hydrocarbon (TPH)  
toxicity characteristic leaching procedure (TCLP)  
unified soil classification system (USCS)  
volatile organic constituent (VOC)

## Executive Summary

The Bloomfield Refinery, which is located in the Four Corners Area of New Mexico, has been in operation since the late 1950s. Past inspections by State and federal environmental inspectors have identified locations where releases to the environment may have occurred. These locations are generally referred to as Solid Waste Management Units (SWMUs) or Areas of Concern (AOCs).

Pursuant to the terms and conditions of an Order issued on July 27, 2007 by the New Mexico Environment Department (NMED) to San Juan Refining Company and Giant Industries Arizona, Inc. for the Bloomfield Refinery, this environmental site investigation was completed for AOC No. 19 (Seep North of MW-45), AOC No. 20 (Seep North of MW-46), and AOC No. 21 (Seep North of MW-47), which are designated as Group 6.

The Order requires that San Juan Refining Company and Giant Industries Arizona, Inc. determine and evaluate the presence, nature, and extent of historical releases of contaminants at the aforementioned AOCs. The investigation activities included collection and analysis of soil samples for potential site-related constituents beginning on August 15, 2011 and continuing through August 17, 2011. The samples were collected from 6 soil borings, which were completed using a hand auger, and from an additional 34 surface sample locations. A total of 45 soil samples (excluding additional quality assurance samples) were collected for analysis of potential site-related constituents (e.g., volatile and semi-volatile organics, total petroleum hydrocarbons, and metals).

Three metals (arsenic, chromium, and cobalt) and diesel range organics were detected in soil samples at concentrations exceeding their respective residential screening levels. Most of the soil samples with constituents detected at concentrations above the screening levels (excluding cobalt) are located at AOC No. 19; cobalt was reported at concentrations above the screening level at all three AOCs. A single sample at both AOC No. 20 and No. 21 had detections of diesel range organics at concentrations above the screening level.

Future studies are planned to evaluate background concentrations at the refinery and after site-specific background concentrations have been established for metals in soils, the reported results should be reevaluated to determine if any of the detected results are representative of background conditions. Additional sampling and analysis should be conducted to determine the

site-specific speciation of chromium in soils near the seeps. Subsequent to the comparison to background and determination of the speciation of chromium, a risk assessment could be completed to provide a more definitive evaluation of the risks posed by the presence of the constituents in area of the seeps.

## **Section 1**

### **Introduction**

The Bloomfield Refinery is located immediately south of Bloomfield, New Mexico in San Juan County (Figure 1). The physical address is #50 Road 4990, Bloomfield, New Mexico 87413. The Bloomfield Refinery is located on approximately 263 acres. Bordering the facility is a combination of federal and private properties. Public property managed by the Bureau of Land Management lies to the south. The majority of undeveloped land in the vicinity of the facility is used extensively for oil and gas production and, in some instances, grazing. U.S. Highway 550 is located approximately one-half mile west of the facility. The topography of the main portion of the site is generally flat with steep bluffs to the north where the San Juan River intersects Tertiary terrace deposits.

The Bloomfield Refinery is currently owned by San Juan Refining Company, a New Mexico corporation, and operated by Western Refining Southwest, Inc. formerly known as Giant Industries Arizona, Inc., an Arizona corporation. The Bloomfield Refinery has an approximate refining capacity of 18,000 barrels per day. Various process units operated at the facility, included crude distillation, reforming, fluidized catalytic cracking, sulfur recovery, mercox treater, catalytic polymerization, and diesel hydrotreating. Products produced at the refinery included gasoline, diesel fuels, jet fuels, kerosene, propane, butane, naphtha, residual fuel, fuel oils, and liquefied petroleum gas (LPG).

On July 27, 2007, the New Mexico Environment Department (NMED) issued an Order to San Juan Refining Company and Giant Industries Arizona, Inc. ("Western") requiring investigation and corrective action at the Bloomfield Refinery. This Investigation Report has been prepared for the AOCs designated as Group 6 in the Order. Group 6 includes AOC No. 19 (Seep North of MW-45), AOC No. 20 (Seep North of MW-46), and AOC No. 21 (Seep North of MW-47).

The area of investigation that is the subject of this report is shown on Figure 2. The Group 6 AOCs are located north of the refinery complex, along the bluff that runs along the south side of the San Juan River.

The purpose of the site investigation is to determine and evaluate the presence, nature, and extent of releases of contaminants in accordance with 20.4.1.500 New Mexico Administrative Code (NMAC) incorporating 40 Code of Federal Regulations (CFR) Section 264.101. The

investigation activities were conducted in accordance with Section IV of the Order and focused on soil as that is the environmental media in these areas that may potentially contain contaminants. Groundwater presence in the area has been significantly reduced by the installation of a slurry wall along the top of bluff in April 2005. The investigation was completed pursuant to the Investigation Work Plan dated December 2009 (revised February 2011), which was approved by the NMED on March 1, 2011.

The samples of soil were analyzed for volatile and semi-volatile organic constituents, petroleum hydrocarbons, and metals. The results of these analyses are compared to applicable State or federal cleanup and screening levels as specified in Section VII. of the Order and as directed by NMED.

## Section 2 Background

This section presents background information for Group 6, including a review of historical waste management activities for each location to identify the following:

- Type and characteristics of waste and contaminants handled in the subject AOCs;
- Known and possible sources of impacts;
- History of releases; and
- Known extent of impacts prior to the current investigation.

From 2001 to 2002, the Hammond Irrigation Ditch was lined with concrete. The ditch extends along the Facility's northern and western boundaries between the refinery and the San Juan River. The water present in the ditch acted as a hydraulic barrier prior to installation of the concrete liner, confining the bulk of the impacts to within the refinery boundaries. A recovery system with a French drain was installed to preserve the integrity of the concrete liner.

Between August and December 2004, petroleum hydrocarbons and water were observed to be seeping out of seven areas on the north side of the refinery near monitoring wells MW-45, MW-46, and MW-47. During inspections of these areas, hydrocarbon was observed to have stained the soil. Figure 2 shows the locations of these seeps.

Four seeps (Nos. 1, 2, 5, and 9) are located within two small drainage tributaries to the San Juan River at a location north of monitoring well MW-45. These seeps are referred to as AOC No. 19 Seep North of MW-45. Two seeps (Seep Nos. 6 and 7) were discovered along the bluff north of monitoring well MW-46 and are referred to as AOC No. 20 Seep North of MW-46. The last seep (No. 8) is located along the bluff above the San Juan River north of monitoring well MW-47 and is referred to as AOC No. 21 Seep North of MW-47.

Nine catchment basins were installed as temporary measures when the seeps were first identified in 2004. A liner was installed at each catchment basin with the exception of Seep No. 4. Initially, the water accumulating in the catchment basins was pumped to Tank 37 and then to the Refinery's wastewater treatment system. The basins were constructed by placing high density polyethylene (HDPE) on top of the exposed bedrock after the visibly impacted soils had been removed and earthen berms formed. The design of the catchment basins also allows the accumulation of rain water in addition to groundwater discharge.

The North Boundary Barrier Wall (slurry wall) was designed and installed along the top of bluff by April 2005 as a permanent control measure to prevent further potential discharges of groundwater to surface water and the generation of seeps along the bluff (Figure 2).

Subsequent to the installation of the slurry wall and with the continued recovery of groundwater from the Hammond Ditch French drain, the discharge from the seeps was significantly reduced.

Only Seeps No. 1, 6, and 9 currently have any actual discharge of groundwater. Seep No. 1 has the most flow at the rate of a trickle. Seep No. 6 has about two thirds the flow of Seep No. 1 with Seep No. 9 measured by drips with no continuous flow. Seep Nos. 2, 5, 7, and 8 did have historic flows, but groundwater discharge has ceased subsequent to the installation of the slurry wall.

Water samples were collected from the seeps in 2007. The samples were collected on three occasions (October 9, 2007, November 8, 2007, and December 10, 2007) from the seeps, which had an active discharge of groundwater. The samples were collected from Seep Nos. 1, 6, 7, 8, and 9 shown on Figure 3 and were identified on the chain-of-custodies and analytical reports as Outfall #1, Outfall #6, Outfall #7, Outfall #8, and Outfall #9. The analytical results are summarized in Table 1. All of the analyses were non-detect for BTEX (benzene, ethylbenzene, toluene, and xylenes) constituents, with one exception. The October sample from Seep No. 1 (Outfall #1) indicated very low concentrations of benzene and xylenes at 0.0029 mg/l and 0.0039 mg/l, respectively. Both subsequent samples from this location were non-detect for all BTEX constituents. All of the detected results and non-detect quantitation limits for the 2007 analyses are less than the potentially applicable groundwater and surface water quality standards shown in Table 1.

Water samples have been collected from the seeps on a semiannual basis starting in 2008. As part of the Facility-Wide Groundwater Monitoring Plan, Western checks all of the identified seep locations (Nos. 1 through 9) on a bi-weekly basis to determine if there is evidence of an active groundwater discharge. The presence of groundwater discharges is recorded and water samples are collected during the semiannual sampling event, when water is present, from all seeps that have indicated an active groundwater discharge within the past six months. The analytical results from these seep samples are summarized in Table 1. No semi-volatile organic compounds were detected in any of the seep samples and, therefore, are not included in the summary table. All analytical results are included in the Groundwater Remediation and Monitoring Report, which is submitted to the NMED in April of each year.

Results of the historical analyses of seep samples show that chloride and sulfate are detected at concentrations above the New Mexico Water Quality Control Commission Standards for Groundwater. All of the 2008 and 2009 analyses were non-detect for BTEX (Western Refining Southwest, Inc., 2009). Methyl tert-butyl ether (MTBE) was sporadically reported in some seep samples at concentrations ranging from less than 0.001 to 0.048 mg/L.



## **Section 3**

### **Scope of Activities**

#### **3.1 Soil Boring Installation and Sample Collection**

Pursuant to Section IV of the Order and the approved Group 6 Investigation Work Plan, an investigation of soils was conducted to determine and evaluate the presence, nature, extent, fate, and transport of contaminants. To accomplish this objective, surface soil samples were collected and soil borings were installed at AOCs 19, 20, and 21 (Figure 3). No groundwater samples were collected as the water from the seeps is already collected and analyzed on a semiannual basis as part of the Facility-Wide Groundwater Monitoring Plan. As outlined in the work plan, there is the potential for constituents to have been released to soils at known locations at AOCs 19, 20, and 21 and therefore a judgmental sampling design was implemented.

Where the HDPE liners are present within the catchment basins, surface soil samples were collected as discussed below and in Section 4.3. Hand augured soil borings were installed at select locations along the arroyos (drainage pathways) and samples were collected from the following intervals where refusal did not prevent sample collection to these depths:

- 0-0.5';
- 1.5-2.5'; and
- Total depth of the boring (if different from the two designated sampling intervals).

Six soil borings were completed at locations in the middle of the drainage pathways and at locations just downstream from the berms of the catchment basins. Thirty four surface soil samples (0-0.5' only) were collected at the seep faces, on top of the HDPE liners, at select locations along the bottom of the drainage pathways, and along the margins of the drainage pathways, which occur on the banks of the drainage pathways and at the edge of the catchment basin liners.

Soil samples were collected at AOC No. 19 from the following locations (Figure 4):

- Three surface soil samples [AOC 19-1 (0-0.5'), AOC 19-2 (0-0.5'), and AOC 19-3 (0-0.5')] were collected near Seep No. 1, with one of the samples collected from the seep face and the other two samples collected along the margins of the drainage pathway;
- Near Seep No.2, a total of three surface soil samples [AOC 19-4 (0-0.5'), AOC 19-5 (0-0.5'), and AOC 19-6 (0-0.5')] were collected with one collected from the seep face and the other two collected from along the margins of the drainage pathway;

- Three surface soil samples [AOC 19-20 (0-0.5'), AOC 19-22 (0-0.5'), and AOC 19-23 (0-0.5')] were collected near Seep No. 5 with one collected from the seep face and two collected along the margins of the drainage pathway;
- Four surface soil samples [AOC 19-14 (0-0.5'), AOC 19-15 (0-0.5'), AOC 19-16 (0-0.5'), and AOC 19-17 (0-0.5')] were collected near Seep No. 9 with one collected from the bottom of the drainage pathway where the seep originated, one sample collected from the center of the lined catchment basin, and two samples collected from along the margins of the drainage pathway;
- One surface soil sample [AOC 19-13 (0-0.5')] was collected at Seep No. 4 at the center of the natural basin where the seep daylighted;
- Two surface soil samples [AOC 19-19 (0-0.5'), and AOC 19-21 (0-0.5')] were collected from along the margins of the drainage pathway between Seep No. 5 and Seep No. 9;
- Four surface soil samples [AOC 19-8 (0-0.5'), AOC 19-10 (0-0.5'), AOC 19-11 (0-0.5'), and AOC 19-12 (0-0.5')] were collected along the margins of the drainage pathway between Seep No. 3 and Seep No. 4; and
- Four soil samples [AOC 19-7 (0-0.5'), AOC 19-7 (1.5-2.0'), AOC 19-9 (0-0.5'), and AOC 19-18 (0-0.5')] were collected from the three soil borings, which are located near the center of the two major drainage pathways that occur at AOC No. 19. The original Scope of Services called for two samples from each boring but refusal was reached at a depth of 0.5' in two of the borings, thus preventing sample collection from the 1.5-2.0' interval.

At AOC No. 20, soil samples were collected at the following locations (Figure 5):

- Four surface samples [AOC 20-1 (0-0.5'), AOC 20-2 (0-0.5'), AOC 20-3 (0-0.5'), and AOC 20-4 (0-0.5')] were collected at Seep No. 6 with one sample collected from the seep face, one soil sample collected from on top of the catchment liner, and two samples collected from along the margins of the drainage pathway;
- Three surface soil samples [AOC 20-9 (0-0.5'), AOC 20-10 (0-0.5'), and AOC 20-11 (0-0.5')] were collected at Seep No. 7, with one sample collected from the top of the catchment liner and two samples collected from along the margins of the drainage pathway;
- Two surface soil samples [AOC 20-8 (0-0.5') and AOC 20-12 (0-0.5')] were collected from along the center line of the drainage pathway that extends from Seep No. 6 toward the San Juan River;
- Two surface soil samples [AOC 20-6 (0-0.5') and AOC 20-7 (0-0.5')] were collected from the margins of the drainage pathway that extends from Seep No. 6 toward the San Juan River; and
- One soil sample [AOC 20-5 (0-0.5')] was collected from a soil boring at the center line of the drainage pathway that extends below Seep No. 6. The original Scope of Services called for two samples from each boring but refusal was reached at a depth of 0.5' in this boring, thus preventing sample collection from the 1.5-2.0' interval.

Soil samples were collected from the following locations at AOC No. 21 (Figure 6):

- Three surface soil samples [AOC 21-2 (0-0.5'), AOC 21-3 (0-0.5'), and AOC 21-4 (0-0.5')] were collected at Seep No. 8, with one collected from the top of the catchment liner and two collected from along the margins of the drainage pathway; and
- Six soil samples [AOC 21-1 (0-0.5'), AOC 21-1 (1.5-2.0'), AOC 21-1 (2.0-2.5'), AOC 21-5 (0-0.5'), AOC 21-5 (1.5-2.0'), and AOC 21-5 (2.0-2.5')] were collected from two soil borings, with one positioned at the discharge point of Seep No. 8 and the other down the drainage pathway just below the catchment berm near Seep No. 8.

### **3.2 Background Information Research**

Documents containing the results of previous investigations and subsequent routine groundwater monitoring data from monitoring wells and the seeps were reviewed to facilitate development of the Investigation Work Plan. The previously collected data provide detailed information on the overall subsurface conditions, including hydrogeology and contaminant distribution within groundwater on a site-wide basis. The data collected under this scope of services will supplement the existing soil and groundwater information and provide specific information regarding contaminant occurrence and distribution within soils near the seeps.

### **3.3 Collection and Management of Investigation Derived Waste**

Soil cuttings (i.e., excess sample material) were contained and will be characterized using methods based on the type of contaminants suspected or encountered. A discrete sample will be collected for waste characterization using a decontaminated stainless-steel spoon because the material is uniform in character and there was only a very small volume of soils generated (i.e., less than one 55-gallon drum), which were collected from within an area with similar types of potential constituents. Additional discussion on management of IDW is presented in Appendix A. The soil sample will be analyzed for leachable RCRA 8 Metals [(Environmental Protection Agency (EPA) method 6010B] using the toxicity characteristic leaching procedure (TCLP), Total Petroleum Hydrocarbons (gasoline, diesel, and motor oil ranges- EPA method 8015B), Polyaromatic Hydrocarbons (EPA method 8310), and BTEX and methyl tertiary butyl ether (EPA method 8260B).

All decontamination water (approximately 30 gallons total) was disposed at the end of each day in the refinery wastewater treatment system upstream of the American Petroleum Institute (API) Separator.

### **3.4 Surveys**

A hand held GPS receiver was used to record the coordinates of each surface soil and soil boring location. These coordinates were recorded on the field boring logs. The soil boring locations were subsequently surveyed by a registered surveyor.

The horizontal coordinates and elevation of each surface sampling location; the surface coordinates and elevation of each soil boring, and the locations of all other pertinent structures was determined by a registered New Mexico professional land surveyor in accordance with the State Plane Coordinate System (NMSA 1978 47-1-49-56 (Repl. Pamp. 1993)). The surveys were conducted in accordance with Sections 500.1 through 500.12 of the Regulations and Rules of the Board of Registration for Professional Engineers and Surveyors Minimum Standards for Surveying in New Mexico. Horizontal positions were measured to the nearest 0.1-ft, and vertical elevations were measured to the nearest 0.01-ft. The survey data is included in Appendix B.

## **Section 4**

### **Field Investigation Results**

This section provides a summary of the surface and subsurface conditions at the refinery, including AOCs No. 19, 20, and 21. A discussion is included on the installation of soil borings, field screening of soils, and collection of soil samples for analysis. Groundwater and surface water conditions are described in the last two subsections.

#### **4.1 Surface Conditions**

Regionally, the surface topography slopes toward the floodplain of the San Juan River, which runs along the northern boundary of the refinery complex. To the south of the refinery, the drainage is to the northwest. North of the refinery, across the San Juan River, surface water flows in a southeasterly direction toward the San Juan River. The active portion of the refinery property, where the process units and storage tanks are located, is generally of low relief with an overall northwest gradient of approximately 0.02 ft/ft. The refinery sits on an alluvial floodplain terrace deposit and there is a steep bluff (approx. drop of 90 feet) at the northern boundary of the refinery where the San Juan River intersects the floodplain terrace, which marks the southern boundary of the floodplain. AOCs 19, 20, and 21 are located on the face of this steep bluff.

There are two locally significant arroyos, one immediately east and another immediately west of the refinery. These arroyos collect most of the surface water flows in the area, thus significantly reducing surface water flows across the refinery. A minor drainage feature is located on the eastern portion of the refinery, where the former Landfill Pond (SWMU No. 9) was located. There are several steep arroyos located to the west along the northern refinery boundary on the steep bluff face that capture local surface water flows and minor groundwater discharges. As shown in the photographs in Appendix C, the seeps associated with AOC Nos. 19, 20 and 21 are located within these western arroyos along the northern refinery boundary.

The refinery complex is bisected by County Rd #4990 (Sullivan Road), which runs east-west. The process units, storage tanks (crude oil and liquid products), wastewater treatment systems, and AOCs 19, 20, and 21 are located north of the county road. The crude oil and product loading racks, LPG storage tanks and loading racks, maintenance buildings/90-day storage area, pipeline offices, transportation truck shop, and the Class I injection well are located south of the county road. There is very little vegetation throughout these areas with most surfaces

composed of concrete, asphalt, or gravel. The area between the refinery and the San Juan River where AOCs 19, 20 and 21 are located does have limited vegetation on steep slopes that do not support dense vegetation.

## **4.2 Subsurface Conditions**

Numerous soil borings and monitoring wells have been completed across the refinery property during previous site investigations and installation of the slurry wall, which runs along the northern and western refinery boundary. Based on the available site-specific and regional subsurface information, the site is underlain by the Quaternary Jackson Lake terrace deposits, which unconformably overlie the Tertiary Nacimiento Formation. The Jackson Lake deposits consist of fine grained sand, silt and clay that grades to coarse sand, gravel and cobble size material closer to the contact with the Nacimiento Formation. The Jackson Lake Formation is over 40 feet thick near the southeast portion of the site and generally thins to the northwest toward the San Juan River. The Nacimiento Formation is primarily composed of fine grained materials (e.g., carbonaceous mudstone/claystone with interbedded sandstones) with a reported local thickness of approximately 570 feet (Groundwater Technology Inc., 1994).

Figures 7 and 8 present cross-sections of the shallow subsurface based on borings logs from on-site monitoring well completions. The boring log for MW-54 on Figure 7 is similar to the logs for MW-45 and MW-46 and depicts the shallow subsurface stratigraphy in the vicinity of AOC Nos. 19, 20 and 21. The surface elevation along the bluff is approximately 5504 feet MSL (based on the surface elevations of monitoring wells MW-45 and MW-46). As shown in Cross-Section A-A' (Figure 7), the top of the Nacimiento Formation occurs at approximately 5495 feet MSL. Due to erosion along the bluff, the elevation of the top of the Nacimiento near the seeps is as much as 60 feet lower.

The uppermost aquifer is under water table conditions and occurs within the sand and gravel deposits of the Jackson Lake Formation. The Nacimiento Formation functions as an aquitard at the site that prevents constituents from migrating to deeper aquifers. The potentiometric surface as measured in August 2009 is presented as Figure 9 and shows the groundwater flowing to the northwest.

Previous investigations have identified and delineated impacts to groundwater from historical site operations. Figure 10 shows the distribution of separate phase hydrocarbon (SPH) in the subsurface based on the apparent thickness of SPH measured in monitoring wells. SPH is not

present within the AOC seeps or within nearby monitoring wells. Dissolved-phase impacts are depicted on Figure 11.

### **4.3 Exploratory Drilling Investigations, Soil Sampling and Boring Abandonment**

This subsection provides a description of surface and subsurface investigations to locate potential impacts to soils. This includes soil field screening results, soil sampling intervals and methods for detection of surface and subsurface impacts in soils.

The soil borings were completed using a hand auger and a general description of the exploratory drilling activities is as follows. The sampling equipment (i.e., hand auger and stainless steel trowel) was decontaminated between each borehole, as described in Appendix A. All soil borings were extended to the depth of refusal, which varied from 0.5 feet to 2.5 feet.

Discrete soil samples were collected for laboratory analysis at the following intervals:

- 0-0.5' (all locations);
- 1.5-2.0' (AOC 19-1, AOC 21-1, and AOC 21-5); and
- 2.0-2.5' (AOC 21-1 and AOC 21-5).

The installation of soil borings and collection of soil samples is discussed below in numerical order for each of the AOCs. A description of the field screening and soil sampling procedures are presented in Appendix A – Field Methods. Copies of the boring logs are provided in Appendix D. In addition to being included on the soil boring logs, the soil vapor (i.e., headspace) screening results are summarized in Table 2. The locations of the soil borings appear on Figures 12 through 19.

#### **4.3.1 AOC 19 – Seep North of MW-45**

##### AOC 19-1

On August 15, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-1. One soil sample was collected from 0 - 0.5 feet. The sample was collected from the discharge point of Seep 1 and was noted as being clayey sand, which was moist to saturated. Field screening with a photoionization detector (PID) indicated potential impacts with an elevated reading of 92 parts per million (ppm).

#### AOC 19-2

On August 15, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-2. One soil sample was collected from 0 - 0.5 feet. The sample was collected near Seep 1 and was noted as being clayey sand, which was damp to dry. Field screening with a PID did not indicate impacts with a reading of 5 ppm and there was no odor.

#### AOC 19-3

One surface soil sample was collected from 0 - 0.5 feet on August 15, 2011 at location AOC 19-3 using a stainless steel trowel. The sample was collected near Seep 1 and was noted as being clayey sand, which was dry to damp. Field screening with a PID did not indicate impacts with a reading of 5 ppm and there was no odor.

#### AOC 19-4

On August 15, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-4. One soil sample was collected from 0 - 0.5 feet. The sample was collected from the historic discharge point of Seep 2 and was noted as being clayey sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 4.2 ppm and there was no odor.

#### AOC 19-5

One surface soil sample was collected from 0 - 0.5 feet on August 15, 2011 at location AOC 19-5 using a stainless steel trowel. The sample was collected near Seep 2 and was noted as being silty sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 3.5 ppm and there was no odor.

#### AOC 19-6

On August 15, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-6. One soil sample was collected from 0 - 0.5 feet. The sample was collected near Seep 2 and was noted as being clayey sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 2 ppm and there was no odor.

#### AOC 19-7

On August 16, 2011 a hand auger was used to collect soil samples at location AOC 19-7 to the depth of refusal, which occurred a depth of 2 feet below ground level (bgl). Soil boring AOC 19-



7 was located along the center line of the drainage pathway below Seep 2 and just above the catchment basin referred to as Seep 3. No odor and low PID readings (2.0 – 2.4 ppm) were observed from 0 to 2 feet bgl in very fine sand and samples were collected from 0 - 0.5 feet and 1.5 - 2.0 feet bgl. Groundwater was not encountered and the boring was plugged on August 16, 2011.

#### AOC 19-8

One surface soil sample was collected from 0 - 0.5 feet on August 15, 2011 at location AOC 19-8 using a stainless steel trowel. The sample was collected along the drainage pathway between catchment basins known as Seep 3 and 4 and was logged as silt, which was dry. Field screening with a PID did not indicate impacts with a reading of 4.8 ppm and there was no odor.

#### AOC 19-9

On August 16, 2011 a hand auger was used to collect soil samples at location AOC 19-9 to the depth of refusal, which occurred a depth of 0.5 feet bgl. Soil boring AOC 19-9 was located along the center line of the drainage pathway between Seep 3 and 4. No odor and a low PID reading (3.2 ppm) were observed from 0 to 0.5 feet bgl in very fine sand and a sample was collected from 0 - 0.5 feet. Groundwater was not encountered and the boring was plugged on August 16, 2011.

#### AOC 19-10

On August 15, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-10. One soil sample was collected from 0 - 0.5 feet. The sample was collected along the drainage pathway between catchment basins known as Seep 3 and 4 and was noted as being fine sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 2.6 ppm and there was no odor.

#### AOC 19-11

One surface soil sample was collected from 0 - 0.5 feet on August 15, 2011 at location AOC 19-11 using a stainless steel trowel. The sample was collected along the drainage pathway just above the catchment basin known as Seep 4 and was logged as silt, which was dry. Field screening with a PID did not indicate impacts with a reading of 3.1 ppm and there was no odor.

#### AOC 19-12

On August 15, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-12. One soil sample was collected from 0 - 0.5 feet. The sample was collected along the drainage pathway just above the catchment basin known as Seep 4 and was noted as being silty fine sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 2.8 ppm and there was no odor.

#### AOC 19-13

One surface soil sample was collected from 0 - 0.5 feet on August 15, 2011 at location AOC 19-13 using a stainless steel trowel. The sample was collected from the center of the catchment basin known as Seep 4 and was logged as silty sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 1.8 ppm and there was no odor.

#### AOC 19-14

On August 16, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-14. One soil sample was collected from 0 - 0.5 feet. The sample was collected near Seep 9 and was logged as silt, which was dry. Field screening with a PID did not indicate impacts with a reading of 2.4 ppm and there was no odor.

#### AOC 19-15

One surface soil sample was collected from 0 - 0.5 feet on August 16, 2011 at location AOC 19-15 using a stainless steel trowel. The sample was collected from center of Seep 9 and was logged as silt, which was dry. Field screening with a PID did not indicate impacts with a reading of 1.2 ppm and there was no odor.

#### AOC 19-16

On August 16, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-16. One soil sample was collected from 0 - 0.5 feet. The sample was collected at Seep 9 and was logged as silty fine sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 4.6 ppm and there was no odor.

#### AOC 19-17

One surface soil sample and a duplicate were collected from 0 - 0.5 feet on August 16, 2011 at location AOC 19-17 using a stainless steel trowel. The sample was collected from the drainage

pathway near the origin of Seep 9 and was logged as sandy clay, which was damp to moist. Field screening with a PID did not indicate impacts with a reading of 3.2 ppm and there was no odor.

#### AOC 19-18

On August 16, 2011 a hand auger was used to collect soil samples at location AOC 19-18 to the depth of refusal, which occurred a depth of 0.5 feet bgl. Soil boring AOC 19-18 was located along the center line of the drainage pathway between Seeps 5 and 9. No odor and a low PID reading (4.2 ppm) were observed from 0 to 0.5 feet bgl in a gravelly, fine to medium grained sand and a sample was collected from 0 - 0.5 feet.

#### AOC 19-19

One surface soil sample was collected from 0 - 0.5 feet on August 16, 2011 at location AOC 19-19 using a stainless steel trowel. The sample was collected from the drainage pathway between Seeps 5 and 9, and was logged as gravelly, very fine sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 2.8 ppm and there was no odor.

#### AOC 19-20

On August 16, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-20. One soil sample was collected from 0 - 0.5 feet. The sample was collected from the drainage pathway at Seep 5 and was logged as gravelly, very fine grained sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 3.2 ppm and there was no odor.

#### AOC 19-21

One surface soil sample was collected from 0 - 0.5 feet on August 16, 2011 at location AOC 19-21 using a stainless steel trowel. The sample was collected between Seeps 5 and 9 and was logged as gravelly, very fine sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 3.6 ppm and there was no odor.

#### AOC 19-22

On August 16, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 19-22. One soil sample was collected from 0 - 0.5 feet. The sample was

collected at Seep 5 and described as gravelly, very fine grained sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 2.1 ppm and there was no odor.

#### AOC 19-23

One surface soil sample was collected from 0 - 0.5 feet on August 16, 2011 at location AOC 19-23 using a stainless steel trowel. The sample was collected at the historic discharge point at Seep 5 and was logged as weathered sandstone, which was dry. Field screening with a PID did not indicate impacts with a reading of 4.9 ppm but there was a faint odor and there was an indication of staining.

### **4.3.2 AOC 20 – Seep North of MW-46**

#### AOC 20-1

On August 17, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 20-1. One soil sample was collected from 0 - 0.5 feet. The sample was collected near the discharge point for Seep 6 and was logged as gravelly, fine grained sand, which was damp. Field screening with a PID did not indicate impacts with a reading of 2.0 ppm and there was no odor.

#### AOC 20-2

One surface soil sample was collected from 0 - 0.5 feet on August 17, 2011 at location AOC 20-2 using a stainless steel trowel. The sample was collected at Seep 6 and was logged as clayey sand, which was damp. Field screening with a PID did not indicate impacts with a reading of 1.3 ppm and there was no odor.

#### AOC 20-3

On August 17, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 20-3. One soil sample and a duplicate were collected from 0 - 0.5 feet. The sample was collected near the center of the containment area for Seep 6 and was logged as clay, which was damp. Field screening with a PID did indicate a potential for impacts with a reading of 33 ppm and there was a faint odor.

#### AOC 20-4

One surface soil sample was collected from 0 - 0.5 feet on August 17, 2011 at location AOC 20-4 using a stainless steel trowel. The sample was collected at Seep 6 and was logged as silt,

which was damp. Field screening with a PID indicated a low potential for impacts with a reading of 9.4 ppm and there was no odor.

#### AOC 20-5

On August 17, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 20-5. This location was designated for a soil boring but refusal was reached at 0.5 feet so the sample was collected with a trowel from 0 - 0.5 feet. The sample was collected along the center of the drainage that flows from Seep 6 to Seep 7. The sample was logged as dry sandy gravel with gravel sizes ranging from ½ inch to 6 inches. Field screening with a PID did not indicate impacts with a reading of 2.4 ppm and there was no odor.

#### AOC 20-6

One surface soil sample was collected from 0 - 0.5 feet on August 17, 2011 at location AOC 20-6 using a stainless steel trowel. The sample was collected along the drainage pathway between Seeps 6 and 7, and was logged as very fine sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 2.4 ppm and there was no odor.

#### AOC 20-7

On August 17, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 20-7. One soil sample was collected from 0 - 0.5 feet. The sample was collected along the drainage that flows from Seep 6 to Seep 7. The sample was logged as dry fine sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 0.7 ppm and there was no odor.

#### AOC 20-8

On August 17, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 20-8. The sample was collected with a trowel from 0 - 0.5 feet. The sample was collected along the center of the drainage that flows from Seep 6 to Seep 7. The sample was logged as dry sandy gravel with gravel sizes ranging from ½ inch to 6 inches. Field screening with a PID did not indicate impacts with a reading of 0.7 ppm and there was no odor.

#### AOC 20-9

One surface soil sample was collected from 0 - 0.5 feet on August 17, 2011 at location AOC 20-9 using a stainless steel trowel. The sample was collected near Seep 7 and was logged as very

fine sand, which was dry. Field screening with a PID did not indicate impacts with a reading of 2.4 ppm and there was no odor.

#### AOC 20-10

On August 17, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 20-10. The sample was collected with a trowel from 0 - 0.5 feet. The sample was collected near Seep 7 and was logged as dry very fine sand. Field screening with a PID did not indicate impacts with a reading of 2.7 ppm and there was no odor.

#### AOC 20-11

One surface soil sample was collected from 0 - 0.5 feet on August 17, 2011 at location AOC 20-11 using a stainless steel trowel. The sample was collected near the center of historic Seep 7 and was logged as dry sandy gravel with gravel sizes ranging from ½ inch to 6 inches. Field screening with a PID did not indicate impacts with a reading of 2.3 ppm and there was no odor.

#### AOC 20-12

On August 17, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 20-12. The sample was collected with a trowel from 0 - 0.5 feet. The sample was collected near Seep 7 and was logged as dry very fine sand. Field screening with a PID did not indicate impacts with a reading of 2.1 ppm and there was no odor.

### **4.3.3 AOC 21 – Seep North of MP-47**

#### AOC 21-1

On August 17, 2011 a hand auger was used to collect soil samples at location AOC 21-1 to the depth of refusal, which occurred a depth of 2.5 feet bgl. Soil boring AOC 21-1 was completed at the historic discharge location of Seep 8. No odor and low PID readings (0.8 – 1.6 ppm) were observed from 0 to 2.5 feet bgl in silty sand. Soil samples were collected from 0 - 0.5 feet, 1.5 - 2.0 feet and 2.0 – 2.5 feet bgl. Groundwater was not encountered and the boring was plugged on August 17, 2011.

#### AOC 21-2

One surface soil sample and a duplicate were collected from 0 - 0.5 feet on August 15, 2011 at location AOC 21-2 using a stainless steel trowel. The sample was collected near Seep 8 and

was logged as dry silty clay. Field screening with a PID indicated a low potential for impacts with a reading of 9.8 ppm and there was no odor.

#### AOC 21-3

On August 15, 2011 a surface soil sample was collected using a stainless steel trowel at location AOC 21-3. The sample was collected from 0 - 0.5 feet. The sample was collected near Seep 8 and was logged as dry silty fine sand. Field screening with a PID did not indicate impacts with a reading of 1.3 ppm and there was no odor.

#### AOC 21-4

One surface soil sample was collected from 0 - 0.5 feet on August 15, 2011 at location AOC 21-4 using a stainless steel trowel. The sample was collected from the center of the catchment basin at Seep 8 and was logged as dry silt. Field screening with a PID did not indicate a potential for impacts with a reading of 5.0 ppm and there was no odor.

#### AOC 21-5

On August 17, 2011 a hand auger was used to collect soil samples at location AOC 21-5 to the depth of refusal, which occurred at a depth of 2.5 feet bgl. Soil boring AOC 21-5 was located along the center of the drainage pathway, downstream of Seep 8. No odor and low PID readings (0.7 ppm) were observed from 0 to 2.5 feet bgl in silty sand. Soil samples were collected from 0 - 0.5 feet, 1.5 - 2.0 feet and 2.0 - 2.5 feet bgl. Groundwater was not encountered and the boring was plugged on August 17, 2011.

### **4.4 Groundwater Conditions**

The uppermost aquifer is under water table conditions and occurs within the sand and gravel deposits of the Jackson Lake Formation. The Nacimiento Formation functions as an aquitard at the site and prevents site related contaminants from migrating to deeper aquifers. The potentiometric surface as measured in August 2009 is presented in Figure 9 and shows the groundwater flowing to the northwest. The potentiometric surface at the site is consistent with the regional gradient in that movement is toward to the San Juan River, which is a location of regional groundwater discharge. The installation of the North Boundary Barrier Wall (i.e., slurry wall) and collection wells/French drain along the western and northern boundary of the refinery controls the flow of groundwater in this area and has basically eliminated the discharge of groundwater from the refinery.

The slurry wall was designed and installed as a permanent control measure to prevent further potential discharges of groundwater to surface water and the generation of seeps along the bluff. Subsequent to the installation of the slurry wall in 2005 and with the continued recovery of groundwater from the Hammond Ditch French drain, the discharge from the seeps was significantly reduced. Only Seeps Nos. 1, 6, and 9 currently have any actual discharge of groundwater. Seep No. 1 has the most flow at the rate of a trickle. Seep No. 6 has about two thirds the flow of Seep No. 1 with Seep No. 9 measured by drips with no continuous flow. Seep Nos. 2, 5, 7, and 8 did have historic flows, but groundwater discharge has ceased subsequent to the installation of the slurry wall.

No separate phase hydrocarbon (SPH) has been observed in the discharge from the seeps since installation of the slurry wall. The areas with the greatest saturated thickness are generally found near and along the Hammond Ditch and on-site surface impoundments (i.e., the current and former raw water ponds). The predominant source of recharge to the shallow aquifer beneath the refinery is recharge from man-made features (e.g., the Hammond Ditch and on-site surface impoundments).

The saturated thickness in the water table aquifer varies from zero feet in the southern and eastern portions of the site to a maximum of approximately eight feet along the northern portion of the refinery, on the up-gradient side of the slurry wall. On the down-gradient side of the slurry wall in the area of AOCs 19, 20, and 21, the saturated thickness has declined significantly since installation of the slurry wall as indicated by the number of dry wells in this area, as shown on Figure 10.

#### **4.5 Surface Water Conditions**

The only local surface water body, excluding on-site surface impounds and the Hammond Irrigation Ditch, is the San Juan River, which flows along the northern most property boundary. There were no accumulations of surface water observed during the site investigation in the areas of AOCs 19, 20, or 21. The discharge from the seeps at AOCs 19, 20, and 21 is currently so low as not to generate an accumulation of surface water but during rainfall events, there could be temporary accumulations of surface water in some of the catchment basins constructed near the seeps.

Regionally, the surface topography slopes toward the floodplain of the San Juan River, and across most of the refinery and to the south of the refinery, the drainage is to the northwest.



There is a steep bluff (approx. drop of 90 feet) at the northern boundary of the refinery where the San Juan River intersects the floodplain terrace, which marks the southern boundary of the floodplain. AOCs 19, 20, and 21 are located on this steep bluff.

There are two locally significant arroyos, one immediately east and another immediately west of the refinery, which collect most of the surface water flows in the area, thus significantly reducing surface water flows across the refinery. A minor drainage feature is located on the eastern portion of the refinery, where the Landfill Pond (SWMU No. 9) was located and there are several steep arroyos along the northern refinery boundary that primarily capture only local surface water flows. There are smaller arroyos located on the bluff face and the seeps are generally near the headwall of the arroyos or along the arroyos.

The average annual rainfall is only approximately 8.6 inches, thus the threat of surface water transport of contaminants as suspended load or dissolved phase is low. The refinery implements a Stormwater Pollution Prevention Plan to ensure that surface waters of the State are not impacted by refinery operations.

## Section 5

### Regulatory Criteria

The applicable screening and cleanup levels are specified in Section VII of the Order issued by NMED on July 2, 2007. The soil cleanup levels are based on a target excess cancer risk of  $10^{-5}$  for carcinogenic contaminants and a target hazard index of 1.0 for noncarcinogenic contaminants. The Order specifies a hierarchy of screening levels, with the screening levels based on NMED guidance taking precedence over EPA's Region VI Human Health Medium Specific Screening Levels with one exception for groundwater. Based on direction received from NMED subsequent to issuance of the Order, EPA's Region VI Human Health Medium Specific Screening Levels have been replaced with EPA Regional Screening Levels dated April 2009. NMED guidance used to establish cleanup levels includes the *Technical Background Document for Development of Soil Screening Levels* (Revision 5.0 dated August 2009) and *Total Petroleum Hydrocarbon (TPH) Screening Guidelines* (dated October 2006).

For non-residential properties (e.g., the Bloomfield Refinery), the soil screening levels must be protective of commercial/industrial workers throughout the upper two feet of surface soils and construction workers throughout the upper ten feet based on NMED criteria. NMED residential soil screening levels are applied to the upper ten feet and soil screening levels for protection of groundwater apply throughout the vadose zone. EPA soil screening levels for direct contact exposure apply to the upper two feet of the vadose zone. To achieve closure as "corrective action complete without controls", the affected media must meet residential screening levels, which are presented in Table 3. Table 4 provides a list of the available NMED and EPA soil screening levels for non-residential properties. While Tables 3 and 4 indicate the various depths to which the individual soil screening levels are applicable, Table 5 discussed below does not include this level of detail.

The aforementioned Tables 3 and 4 have soil screening levels for the soil-to-groundwater pathway that are based on a dilution/attenuation factor (DAF) of 1.0 and 11.25. A review of site conditions indicates that a DAF of 1.0 is overly conservative, thus a site-specific DAF value was calculated. A review of the site-specific conditions at each of the SWMUs and AOCs recently investigated indicates that the conditions at SWMU No. 2 Drum Storage Area North Bone Yard could present a greater potential for constituents to leach from soils to the underlying groundwater because SWMU No. 2 has the shallowest depth to groundwater. There is not a

significant accumulation of groundwater present above the Nacimiento Formation at AOCs 19, 20, or 21. A DAF value of 11.25 was calculated for SWMU No. 2 in the Group No. 2 Investigation Report (dated May 2009, revised March 2010) and although it may be overly conservative for the AOCs investigated under Group 6, the same DAF value of 11.25 is applied at all locations presented in this Investigation Report. The documentation of the calculation of the site-specific DAF value is provided in Appendix E.

The screening levels that are compared to individual soil sample results from AOCs 19, 20, and 21 are presented in Table 5. The screening levels included in Table 5 are based on residential land use and include screening levels to evaluate the potential for constituents to migrate to groundwater using the site-specific DAF of 11.25. The screening levels in Table 5 have not been segregated based on depth of the soil sample as discussed above for Tables 3 and 4. It should also be noted that the screening levels for chromium are conservatively based on the presence of chromium VI; speciation has not been conducted to determine if the detected chromium is chromium III or chromium VI.

A review of the NMED TPH Screening Guidelines (dated 2006) indicates that the TPH screening levels were developed based on screening levels and compositional assumptions developed by the Massachusetts Department of Environmental Protection (MADEP). The screening levels were developed to be protective of both direct contact exposures and leachate generation. The analytical results, as presented in Table 5, are reported for gasoline range organics (C6- C10), diesel range organics (>C10-C28), and motor oil range organics (>C28-C35). The applicable TPH screening levels used for comparison to the individual samples are selected from Table 2a of the NMED guidance. Because the specific type of products (e.g., diesel fuel or jet fuel) that were released in the potential source areas and migrated to AOCs 19, 20, or 21 is not known, "unknown oil" was selected from Table 2a for comparison to the diesel range analytical results. There are no screening levels for comparison to results reported in the gasoline range.

The analyses for motor oil range organics only report results for >C28 to C35. Since the motor oil range analytical results only include hydrocarbons greater than C28, it is not appropriate to compare the results against screening levels for product types that have lower hydrocarbon ranges (e.g., diesel fuel – 60% C11-C22 aromatics and 40% C9-C18 aliphatics). The only product type in the NMED guidance that contains the >C28-C35 carbon range is "waste oil",

which includes C19-C36. Therefore, the motor oil range soil analytical results are compared to the “waste oil” soil screening levels.

The DAF screening levels for diesel and motor range organics in Table 5 are conservatively based on residential land use. The DAF screening levels have not been adjusted for the site-specific DAF value of 11.25 because the default screening levels that NMED based on the MADEP screening levels already incorporate a minor adjustment for the DAF component.

To evaluate the potential for “direct contact” type exposures (e.g., dermal contact, ingestion, and inhalation of particulates and volatiles) to TPH in surface soils, a screening level was developed for the “direct contact” pathways. This screening level was developed for C11-C22 aromatics as it has one of the highest toxicities of any of the TPH fractions used by NMED to calculate screening levels and is used to compare to “unknown oil.” The calculation of the screening level for C11-C22 aromatics is documented in Appendix F. The calculation uses Equation 1 (Combined Exposures to Noncarcinogenic Contaminants in Soil; Residential Scenario) and all provided default exposure values from NMED’s August 2009 *Technical Background Document for Development of Soil Screening Levels, Revision 5.0*. The toxicity values are taken from MADEP’s Massachusetts Contingency Plan Standards spreadsheets, 2009 (<http://www.mass.gov/dep/cleanup/laws/standard.htm>). This screening level of 1,830 mg/kg is shown in the “residential soil screening level” column of Table 5 and is used for comparison to the diesel range analytical results. This calculation was previously approved by NMED in the Group 2 Investigation Report (revised March 2010). A separate “direct contact” screening level was not developed for motor oil because all detected values are already less than the more conservative DAF screening level.

Some of the individual constituents reported by the laboratory do not have screening levels but were all non-detect except 2-hexanone, 4-methyl-2-pentanone, and 4-isopropyltoluene. These constituents were detected at low concentrations (e.g., a maximum of 0.00584 mg/kg) in only a few samples.

## Section 6

### Site Impacts

This section discusses the chemical analyses performed and presents the analytical results that were obtained through the analysis of soil samples, which were collected at the Group 6 AOCs. The analytical results for soils are presented and compared to applicable screening levels, as described in Section 5.0.

#### 6.1 Soil Sampling Chemical Analytical Results

Soil samples were analyzed by Hall Environmental Analysis Laboratory in Albuquerque, New Mexico using the following methods for organic constituents:

- SW-846 Method 8260 volatile organic compounds;
- SW-846 Method 8270 semi-volatile organic compounds; and
- SW-846 Method 8015B gasoline, diesel, and motor oil range petroleum hydrocarbons.

Soil samples were analyzed for the following metals using the indicated analytical methods.

| 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       |
| Cobalt    | SW-846 method 6010/6020       |
| Cyanide   | SW-846 method 335.3/335.2 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       |

The analytical results for AOCs 19, 20, and 21 are summarized in Table 5. The individual results that exceed the applicable screening levels are indicated using a bold font and/or are highlighted, as explained in the table footnotes. Maps showing the distribution of constituents detected in soils above screening levels are included as Figures 12 – 19. The concentrations

shown on Figures 12 – 19 that exceed the screening levels in Table 5 are underlined on the figures (i.e., concentrations above residential and/or DAF (11.25) screening levels). The laboratory analytical reports are included in Appendix G and the data validation of the results, which includes the analytical results for the associated QA/QC samples, is included in Appendix H. The constituents that have concentrations in soils above screening levels are discussed below.

#### AOC No. 19 – Seep North of MW-45

Arsenic was detected in nine samples [AOC 19-2 (0-0.5'), AOC 19-3 (0-0.5'), AOC 19-8 (0-0.5'), AOC 19-9 (0-0.5'), AOC 19-13 (0-0.5'), AOC 19-16 (0-0.5'), AOC 19-17 (0-0.5'), AOC 19-18 (0-0.5'), and AOC 19-23 (0-0.5')] at concentrations above the DAF screening level (0.148 mg/kg) and all of these samples except AOC 19-16 (0-0.5') also have concentrations above the residential screening level (3.59 mg/kg). The detected concentrations range from 2.9 mg/kg to 24 mg/kg. The arsenic analytical results are shown on Figure 12.

Chromium (reported as total chromium) was detected in five soil samples [AOC 19-13 (0-0.5'), AOC 19-14 (0-0.5'), AOC 19-16 (0-0.5'), AOC 19-17 (0-0.5'), and AOC 19-18 (0-0.5')] at a concentration above the chromium VI DAF screening level of 23.7 mg/kg. Chromium was detected in one soil sample [AOC 19-14 (0-0.5')] at a concentration of 600 mg/kg, which exceeds the chromium VI residential soil screening level of 219 mg/kg. Speciation of the detected chromium results was not performed to determine if the chromium is present as chromium III or chromium VI. The lowest chromium III screening level is a value of 113,000 mg/kg for residential exposure. The detected concentrations ranged from 1.6 mg/kg to 600 mg/kg. The chromium analytical results are shown on Figure 13.

Cobalt was detected in three soil samples [(AOC 19-2 (0-0.5'), AOC 19-12 (0-0.5'), and AOC 19-15 (0-0.5')] at concentrations above the DAF screening level (5.51 mg/kg) but none exceed the residential soil screening levels. The detected concentrations range from 1.6 mg/kg to 11 mg/kg. The cobalt analytical results are shown on Figure 14.

Diesel range organics were detected in four samples [(AOC 19-1 (0-0.5'), AOC 19-13 (0-0.5'), AOC 19-14 (0-0.5'), and AOC 19-23 (0-0.5')] at concentrations above the DAF soil screening level and one of these samples [AOC 19-23 (0-0.5')] had a concentration of 21,000 mg/kg, which exceeds the residential direct contact screening level of 1,830 mg/kg. The detected concentrations range from 16 mg/kg to 21,000 mg/kg. The diesel range organics analytical results are shown on Figure 15.

#### AOC No. 20 – Seep North of MW-46

Cobalt was detected in one soil sample [(AOC 20-3 (0-0.5'))] at a concentration (8.9 mg/kg) above the DAF screening level (5.51 mg/kg) but did not exceed the residential soil screening level of 23 mg/kg. The cobalt analytical results for AOC No. 20 are shown on Figure 16.

Diesel range organics were detected in one sample [(AOC 20-4 (0-0.5'))] at a concentration (490 mg/kg) above the DAF soil screening level of 200 mg/kg. None of the analytical results exceeded the residential direct contact screening level of 1,830 mg/kg. The diesel range organics analytical results for AOC No. 20 are shown on Figure 17.

#### AOC No. 21 – Seep North of MW-47

Cobalt was detected in eight soil samples [(AOC 21-1 (0-0.5'), AOC 21-1 (1.5-2'), AOC 21-2 (0-0.5'), AOC 21-3 (0-0.5'), AOC 21-4 (0-0.5'), AOC 21-5 (0-0.5'), AOC 21-5 (1.5-2'), and AOC 21-5 (2.0-2.5')] at concentrations above the DAF screening level (5.51 mg/kg) but none exceed the residential soil screening level of 23 mg/kg. The detected concentrations range from 5.5 mg/kg to 9.1 mg/kg. The cobalt analytical results for AOC No. 21 are shown on Figure 18.

Diesel range organics were detected in one sample [(AOC 21-2 (0-0.5'))] at a concentration (710 mg/kg) above the DAF soil screening level of 200 mg/kg. None of the analytical results exceeded the residential direct contact screening level of 1,830 mg/kg. The diesel range organics analytical results for AOC No. 21 are shown on Figure 19.

## Section 7

# Conclusions and Recommendations

This section summarizes and provides an evaluation of the potential impacts as shown in field screening data and analytical data. An investigation of soils was conducted at Group 6 (AOCs 19, 20, and 21) to assess and evaluate the presence, nature, extent, fate, and transport of contaminants. To accomplish this objective, soil samples were collected and analyzed for potential site-related constituents.

### 7.1 Conclusions

#### AOC No. 19 – Seep North of MW-45

Three metals (arsenic, cobalt, and chromium) and diesel range organics were detected in soil samples at concentrations exceeding their respective screening levels. Arsenic was detected in nine samples at concentrations above the DAF screening level, with eight of these samples having concentrations above the residential screening level. Cobalt was detected in generally low concentrations with a maximum detected value of 11 mg/kg. As shown in Table 5, the samples with concentrations of cobalt above the screening levels do not have concentrations of organic constituents above screening levels. Chromium was detected in five samples at concentrations above the DAF screening level with only one being above the residential screening level. Arsenic, chromium, and cobalt are naturally occurring constituents in many soils and it is possible that these constituents are naturally occurring in soils at the Bloomfield Refinery. The analyses for chromium were for total chromium, which are compared to the chromium VI screening levels. It may be overly conservative to compare to the total chromium results to the chromium VI screening levels, but additional analyses will be required to speciate the chromium.

The only analytical results for organic constituents (i.e., diesel range organics) above the DAF screening level occur at three locations with a maximum value of 770 mg/kg. All of the detections of diesel range organics are below the residential screening level.

#### AOC No. 20 – Seep North of MW-46

Only cobalt and diesel range organics were detected at concentrations above the applicable screening levels. Cobalt was detected in only one soil sample at a concentration (8.9 mg/kg)



above the DAF screening level (5.51 mg/kg) but did not exceed the residential soil screening level of 23 mg/kg. As noted above, cobalt may be a naturally occurring constituent. Diesel range organics were detected in only one sample at a concentration (490 mg/kg) above the DAF soil screening level of 200 mg/kg. None of the diesel range organic analytical results exceeded the residential direct contact screening level of 1,830 mg/kg.

#### AOC No. 21 – Seep North of MW-47

Similarly to AOC No. 20, only cobalt and diesel range organics were detected at concentrations above the applicable screening levels. Cobalt was detected in eight soil samples at concentrations above the DAF screening level (5.51 mg/kg) but none exceed the residential soil screening level of 23 mg/kg. Diesel range organics were detected in one sample at a concentration (710 mg/kg) above the DAF soil screening level of 200 mg/kg. None of the diesel range organic analytical results exceed the residential direct contact screening level of 1,830 mg/kg.

#### Cumulative Risk Evaluation

A cumulative risk evaluation for soils is presented in Table 6. This was conducted by taking the maximum reported concentration of each detected constituent from all three AOCs and dividing by the residential screening level and non-residential screening levels as shown in the equations below. These calculations are separated for carcinogenic and non-carcinogenic constituents. The cumulative carcinogenic risk is  $6.703 \times 10^{-5}$  assuming residential land use and  $1.36 \times 10^{-5}$  for non-residential land use. The hazard index for residential land use is 3.557 and for non-residential land use is 1.652.

$$\text{Site Risk} = \left( \frac{\text{conc}_x}{\text{SSL}_x} + \frac{\text{conc}_y}{\text{SSL}_y} + \frac{\text{conc}_z}{\text{SSL}_z} + \dots + \frac{\text{conc}_i}{\text{SSL}_i} \right) \times 10^{-5}$$

$$\text{Site Hazard Index (HI)} = \left( \frac{\text{conc}_x}{\text{SSL}_x} + \frac{\text{conc}_y}{\text{SSL}_y} + \frac{\text{conc}_z}{\text{SSL}_z} + \dots + \frac{\text{conc}_i}{\text{SSL}_i} \right) \times 1$$

## **7.2 Recommendations**

Western plans to conduct an investigation of background locations and will attempt to establish site-specific background values for various naturally occurring constituents (e.g., arsenic and

cobalt in soils). After background values have been established, site data collected for AOCs 19, 20 and 21 should be reassessed to determine if some analytical results are actually representative of background conditions and not site-related impacts.

Additional soil samples should be collected at the locations with the highest concentrations of chromium at AOC 19 [e.g., AOC19-13 (0-0.5'), AOC 19-14 (0-0.5'), and AOC 19-18 (0-0.5')]. The samples should be analyzed for chromium VI to determine if the chromium detected in soils at AOC 19 is present as chromium III or chromium VI. If the chromium is determined to be chromium III instead of chromium VI, then the screening levels in Table 5 will need to be revised along with any associated figures.

After site-specific background values have been established and the location-specific speciation of chromium has been determined, then preparation of a risk assessment for the Group 6 AOCs may be appropriate. The site concentrations have been compared to residential screening levels in this investigation report; however, the physical nature of the seep areas may limit future use of this area. An evaluation of potential exposures in a site-specific risk assessment could offer a more appropriate evaluation of actual risks based on more likely future use of the area at AOCs 19, 20, and 21.

## **Section 8**

### **References**

Groundwater Technology Inc., 1994, RCRA Facility Investigation/Corrective Measures Study Report Bloomfield Refining Company #50 County Road 4990 Bloomfield, New Mexico, p.51.

Western Refining Southwest, Inc., 2009, Groundwater Remediation and Monitoring Annual Report, Bloomfield Refinery Bloomfield New Mexico, p.37.

Table 1  
Group 6 Historical Seep Discharge Analyses  
Western Refining Southwest - Bloomfield Refinery

| Seep Location | Date   | EPA 8260B          |                   |                    |                   |                    | EPA 300.0          |                    |                   |                   |                   |                      |                   | SM 2320B        |                   |
|---------------|--------|--------------------|-------------------|--------------------|-------------------|--------------------|--------------------|--------------------|-------------------|-------------------|-------------------|----------------------|-------------------|-----------------|-------------------|
|               |        | Benzene<br>(mg/L)  | Toluene<br>(mg/L) | EthylBen<br>(mg/L) | Xylene<br>(mg/L)  | MTBE<br>(mg/L)     | Fluoride<br>(mg/L) | Chloride<br>(mg/L) | Bromide<br>(mg/L) | Nitrite<br>(mg/L) | Nitrate<br>(mg/L) | Phosphorus<br>(mg/L) | Sulfate<br>(mg/L) | CO2<br>(mg/L)   | Arsenic<br>(mg/L) |
|               |        | 0.005 <sup>2</sup> | 0.75 <sup>1</sup> | 0.7 <sup>2</sup>   | 0.62 <sup>1</sup> | 0.012 <sup>3</sup> | 1.6 <sup>1</sup>   | 250 <sup>1</sup>   | NE                | 1 <sup>2</sup>    | 10 <sup>2</sup>   | NE                   | 600 <sup>1</sup>  | NE              | NE                |
| 1             | Aug-09 | <0.001             | <0.001            | <0.001             | <0.003            | 0.003              | 0.29               | 390                | NA                | <2.0              | <0.10             | <0.50                | 1500              | 200             | 2                 |
|               | Apr-09 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>    | NS <sup>1</sup>    | NA                | NS <sup>1</sup>   | NS <sup>1</sup>   | NS <sup>1</sup>      | NS <sup>1</sup>   | NS <sup>1</sup> | 1                 |
|               | Aug-08 | <0.001             | <0.001            | <0.001             | <0.003            | 0.042              | 0.35               | 370                | NA                | <1.0              | <0.10             | <0.50                | 1500              | 250             | 2                 |
|               | Apr-08 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NA                | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Dec-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Nov-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Oct-07 | <0.001             | 0.0029            | <0.001             | 0.0039            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
| 3             | Aug-09 | <0.001             | <0.001            | <0.001             | <0.003            | <0.0015            | 0.22               | 1400               | NA                | <2.0              | <2.0              | <10                  | 6800              | 120             | 1                 |
|               | Apr-09 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>    | NS <sup>1</sup>    | NA                | NS <sup>1</sup>   | NS <sup>1</sup>   | NS <sup>1</sup>      | NS <sup>1</sup>   | NS <sup>1</sup> | 1                 |
|               | Aug-08 | <0.001             | <0.001            | <0.001             | <0.003            | <0.015             | 0.8                | 370                | NA                | <1.0              | <0.10             | <0.50                | 2500              | 160             | 1                 |
|               | Apr-08 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NA                | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
| 6             | Aug-09 | <0.001             | <0.001            | <0.001             | <0.003            | <0.0015            | 0.58               | 4800               | NA                | <2.0              | <2.0              | <10                  | 1500              | 150             | 1                 |
|               | Apr-09 | <0.001             | <0.001            | <0.001             | <0.002            | 0.014              | <0.50              | 2900               | 1.6               | <2.0              |                   | <0.50                | 1000              | 420             | 4                 |
|               | Aug-08 | <0.001             | <0.001            | <0.001             | <0.003            | 0.006              | 0.47               | 2500               | NA                | <1.0              | <0.10             | <0.50                | 960               | 370             | 3                 |
|               | Apr-08 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NA                | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Dec-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Nov-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Oct-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
| 7             | Aug-09 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>    | NS <sup>1</sup>    | NA                | NS <sup>1</sup>   | NS <sup>1</sup>   | NS <sup>1</sup>      | NS <sup>1</sup>   | NS <sup>1</sup> | 1                 |
|               | Apr-09 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>    | NS <sup>1</sup>    | NA                | NS <sup>1</sup>   | NS <sup>1</sup>   | NS <sup>1</sup>      | NS <sup>1</sup>   | NS <sup>1</sup> | 1                 |
|               | Aug-08 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>    | NS <sup>1</sup>    | NA                | NS <sup>1</sup>   | NS <sup>1</sup>   | NS <sup>1</sup>      | NS <sup>1</sup>   | NS <sup>1</sup> | 1                 |
|               | Apr-08 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NA                | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Dec-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Nov-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Oct-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
| 8             | Aug-09 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>    | NS <sup>1</sup>    | NA                | NS <sup>1</sup>   | NS <sup>1</sup>   | NS <sup>1</sup>      | NS <sup>1</sup>   | NS <sup>1</sup> | 1                 |
|               | Apr-09 | <0.001             | <0.001            | <0.001             | <0.002            | <0.001             | 0.33               | 650                | 1.7               | <1.0              |                   | <0.50                | 2200              | 190             | 2                 |
|               | Aug-08 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>    | NS <sup>1</sup>    | NA                | NS <sup>1</sup>   | NS <sup>1</sup>   | NS <sup>1</sup>      | NS <sup>1</sup>   | NS <sup>1</sup> | 1                 |
|               | Apr-08 | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NA                | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Dec-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Nov-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Oct-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
| 9             | Aug-09 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>    | NS <sup>1</sup>    | NA                | NS <sup>1</sup>   | NS <sup>1</sup>   | NS <sup>1</sup>      | NS <sup>1</sup>   | NS <sup>1</sup> | 1                 |
|               | Apr-09 | <0.001             | <0.001            | <0.001             | <0.002            | 0.048              | 0.35               | 620                | 1.0               | <1.0              |                   | <0.50                | 1500              | 420             | 4                 |
|               | Aug-08 | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>   | NS <sup>1</sup>    | NS <sup>1</sup>    | NS <sup>1</sup>    | NA                | NS <sup>1</sup>   | NS <sup>1</sup>   | NS <sup>1</sup>      | NS <sup>1</sup>   | NS <sup>1</sup> | 1                 |
|               | Apr-08 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NA                | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Dec-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Nov-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |
|               | Oct-07 | <0.001             | <0.001            | <0.001             | <0.002            | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>    | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>   | NS <sup>2</sup>      | NS <sup>2</sup>   | NS <sup>2</sup> | 1                 |

NE = not established  
NS<sup>1</sup> = Seep location is Dry or Not Enough Water to Sample- No Sample  
NS<sup>2</sup> = Not Sampled - Sample was taken before implementation of Facility-Wide Monitoring Plan  
NA = Constituent analysis not requested  
1- WQCC 20NMAC 6.2.3103 = New Mexico Water Quality Control Commission Standard for Groundwater of 10,000 ug/l  
2 - 40 CFR141.61 = Safe Driking Water Act Maximum Contaminant Level  
3 - EPA Regional Screening Level - Tap Water  
No active discharge at Seeps 2, 4 and 5 has been present to sample since implementation of the Facility-Wide Monitoring  
mg/L = milligram per liter  

0.048

 - bolded and shaded value exceeds screening level

**TABLE 2**  
**Group 6 Soil Boring Samples - Vapor Screening Results**  
**Bloomfield Refinery - Bloomfield, New Mexico**

| AOC 19                           | OVM (ppm) | AOC 20    | OVM (ppm) | AOC 21   | OVM (ppm) |
|----------------------------------|-----------|-----------|-----------|----------|-----------|
| Sample Interval Depth - 0-0.5'   |           |           |           |          |           |
| AOC 19-1                         | 92        | AOC 20-1  | 2         | AOC 21-1 | 1.5       |
| AOC 19-2                         | 5         | AOC 20-2  | 1.3       | AOC 21-2 | 9.8       |
| AOC 19-3                         | 5         | AOC 20-3  | 33        | AOC 21-3 | 1.3       |
| AOC 19-4                         | 4.2       | AOC 20-4  | 9.4       | AOC 21-4 | 5         |
| AOC 19-5                         | 3.5       | AOC 20-5  | 2.4       | AOC 21-5 | 0.7       |
| AOC 19-6                         | 2         | AOC 20-6  | 2.4       |          |           |
| AOC 19-7                         | 2.3       | AOC 20-7  | 0.7       |          |           |
| AOC 19-8                         | 4.8       | AOC 20-8  | 0.7       |          |           |
| AOC 19-9                         | 3.2       | AOC 20-9  | 2.4       |          |           |
| AOC 19-10                        | 2.6       | AOC 20-10 | 2.7       |          |           |
| AOC 19-11                        | 3.1       | AOC 20-11 | 2.3       |          |           |
| AOC 19-12                        | 2.8       | AOC 20-12 | 2.1       |          |           |
| AOC 19-13                        | 1.8       |           |           |          |           |
| AOC 19-14                        | 2.4       |           |           |          |           |
| AOC 19-15                        | 1.2       |           |           |          |           |
| AOC 19-16                        | 4.6       |           |           |          |           |
| AOC 19-17                        | 3.2       |           |           |          |           |
| AOC 19-18                        | 4.2       |           |           |          |           |
| AOC 19-19                        | 2.8       |           |           |          |           |
| AOC 19-20                        | 3.2       |           |           |          |           |
| AOC 19-21                        | 3.6       |           |           |          |           |
| AOC 19-22                        | 2.1       |           |           |          |           |
| AOC 19-23                        | 4.9       |           |           |          |           |
| Sample Interval Depth - 0.5-1.5' |           |           |           |          |           |
| AOC 19-7                         | 2         |           |           | AOC 21-1 | 0.8       |
|                                  |           |           |           | AOC 21-5 | 0.7       |
| Sample Interval Depth - 1.5-2'   |           |           |           |          |           |
| AOC 19-7                         | 2.4       |           |           | AOC 21-1 | 1.6       |
|                                  |           |           |           | AOC 21-5 | 0.7       |
| Sample Interval Depth - 2-2.5'   |           |           |           |          |           |
|                                  |           |           |           | AOC 21-1 | 1.1       |
|                                  |           |           |           | AOC 21-5 | 0.7       |

OVM - organic vapor monitor  
ppm - parts per million

**TABLE 3**  
**Residential Soil Screening Levels**  
**Bloomfield Refinery - Bloomfield, New Mexico**

| Analyte                     | NMED                     |          | EPA                      |             | Cross Media Soil-to-Ground Water |                           |                          | Cross Media Soil-to-Ground Water |                                   |                                  | Constituent Detected |
|-----------------------------|--------------------------|----------|--------------------------|-------------|----------------------------------|---------------------------|--------------------------|----------------------------------|-----------------------------------|----------------------------------|----------------------|
|                             | Residential Soil (mg/kg) | Endpoint | Residential Soil (mg/kg) | ResSoil key | DAF1 (mg/kg)                     | GW_Risk-based SSL (mg/kg) | GW_MCL-based SSL (mg/kg) | DAF (11.25) (mg/kg)              | GW_Risk-based SSL (11.25) (mg/kg) | GW_MCL-based SSL (11.25) (mg/kg) |                      |
|                             |                          |          |                          |             |                                  |                           |                          |                                  |                                   |                                  |                      |
| Applicable depth interval   | 0-10'                    |          | 0-2'                     |             | All depths                       |                           |                          | All depths                       |                                   |                                  |                      |
| Acenaphthene                | 3.44E+03                 | ns       | 3.40E+03                 | n           | 2.05E+01                         | 2.70E+01                  | -                        | 2.31E+02                         | 3.04E+02                          | -                                | N                    |
| Acenaphthylene              | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Acetone                     | 6.75E+04                 | n        | 6.10E+04                 | n           | 3.84E+00                         | 4.40E+00                  | -                        | 4.32E+01                         | 4.95E+01                          | -                                | Y                    |
| Aniline                     | -                        | -        | 8.50E+01                 | c**         | -                                | 3.40E-03                  | -                        | -                                | 3.83E-02                          | -                                | N                    |
| Anthracene                  | 1.72E+04                 | ns       | 1.70E+04                 | n           | 3.37E+02                         | 4.50E+02                  | -                        | 3.79E+03                         | 5.06E+03                          | -                                | N                    |
| Antimony                    | 3.13E+01                 | n        | 3.10E+01                 | n           | 6.61E-01                         | 6.60E-01                  | 2.70E-01                 | 7.44E+00                         | 7.43E+00                          | 3.04E+00                         | N                    |
| Arsenic                     | 3.59E+00                 | c        | 3.90E-01                 | c*          | 1.31E-02                         | 1.30E-03                  | 2.90E-01                 | 1.48E-01                         | 1.46E-02                          | 3.26E+00                         | Y                    |
| Azobenzene                  | -                        | -        | 4.90E+00                 | c           | -                                | 5.10E-04                  | -                        | -                                | 5.74E-03                          | -                                | N                    |
| Barium                      | 1.56E+04                 | n        | 1.50E+04                 | n           | 3.01E+02                         | 3.00E+02                  | 8.20E+01                 | 3.39E+03                         | 3.38E+03                          | 9.23E+02                         | Y                    |
| Benz(a)anthracene           | 4.81E+00                 | c        | 1.50E-01                 | c           | 3.20E-01                         | 1.40E-02                  | -                        | 3.59E+00                         | 1.58E-01                          | -                                | N                    |
| Benzene                     | 1.55E+01                 | c        | 1.10E+00                 | c*          | 1.85E-03                         | 2.30E-04                  | 2.80E-03                 | 2.08E-02                         | 2.59E-03                          | 3.15E-02                         | N                    |
| Benzo(a)pyrene              | 4.81E-01                 | c        | 1.50E-02                 | c           | 1.09E-01                         | 4.60E-03                  | 3.10E-01                 | 1.22E+00                         | 5.18E-02                          | 3.49E+00                         | N                    |
| Benzo(b)fluoranthene        | 4.81E+00                 | c        | 1.50E-01                 | c           | 1.11E+00                         | 4.70E-02                  | -                        | 1.25E+01                         | 5.29E-01                          | -                                | N                    |
| Benzo(g,h,i)perylene        | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Benzo(k)fluoranthene        | 4.81E+01                 | c        | 1.50E+00                 | c           | 1.09E+01                         | 4.60E-01                  | -                        | 1.22E+02                         | 5.18E+00                          | -                                | N                    |
| Benzoic acid                | -                        | -        | 2.40E+05                 | nm          | -                                | 3.30E+01                  | -                        | -                                | 3.71E+02                          | -                                | Y                    |
| Benzyl alcohol              | -                        | -        | 3.10E+04                 | n           | -                                | 4.20E+00                  | -                        | -                                | 4.73E+01                          | -                                | N                    |
| Beryllium                   | 1.56E+02                 | n        | 1.60E+02                 | n           | 5.77E+01                         | 5.80E+01                  | 3.20E+00                 | 6.49E+02                         | 6.53E+02                          | 3.60E+01                         | Y                    |
| Bis(2-chloroethoxy)methane  | -                        | -        | 1.80E+02                 | n           | -                                | 2.30E-02                  | -                        | -                                | 2.59E-01                          | -                                | N                    |
| Bis(2-chloroethyl)ether     | 2.56E+00                 | c        | 1.90E-01                 | c           | 2.33E-05                         | 2.70E-06                  | -                        | 2.62E-04                         | 3.04E-05                          | -                                | N                    |
| Bis(2-chloroisopropyl)ether | 9.15E+01                 | c        | -                        | -           | 2.56E-03                         | -                         | -                        | 2.88E-02                         | -                                 | -                                | N                    |
| Bis(2-ethylhexyl)phthalate  | 2.80E+02                 | c        | 3.50E+01                 | c*          | 1.19E+01                         | 1.60E+00                  | 2.00E+00                 | 1.34E+02                         | 1.80E+01                          | 2.25E+01                         | N                    |
| Bromobenzene                | -                        | -        | 9.40E+01                 | n           | -                                | 1.50E-02                  | -                        | -                                | 1.69E-01                          | -                                | N                    |
| Bromodichloromethane        | 5.25E+00                 | c        | 2.80E-01                 | c           | 2.76E-04                         | 3.30E-05                  | -                        | 3.11E-03                         | 3.71E-04                          | -                                | N                    |
| Bromoform                   | -                        | -        | 6.10E+01                 | c*          | -                                | 2.30E-03                  | -                        | -                                | 2.59E-02                          | -                                | N                    |
| Bromomethane                | 2.23E+01                 | n        | 7.90E+00                 | n           | 1.94E-03                         | 2.20E-03                  | -                        | 2.18E-02                         | 2.48E-02                          | -                                | N                    |
| 4-Bromophenyl phenyl ether  | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 2-Butanone (MEK)            | 3.96E+04                 | n        | 2.80E+04                 | ns          | 1.27E+00                         | 1.50E+00                  | -                        | 1.43E+01                         | 1.69E+01                          | -                                | Y                    |
| Butyl benzyl phthalate      | -                        | -        | 2.60E+02                 | c*          | -                                | 6.70E-01                  | -                        | -                                | 7.54E+00                          | -                                | N                    |
| Cadmium                     | 7.79E+01                 | n        | 7.00E+01                 | n           | 1.37E+00                         | 1.40E+00                  | 3.80E-01                 | 1.55E+01                         | 1.58E+01                          | 4.28E+00                         | N                    |
| Carbazole                   | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Carbon disulfide            | 1.94E+03                 | ns       | 6.70E+02                 | ns          | 2.52E-01                         | 2.70E-01                  | -                        | 2.84E+00                         | 3.04E+00                          | -                                | Y                    |
| Carbon tetrachloride        | 4.38E+00                 | c        | 2.50E-01                 | c           | 7.39E-04                         | 7.90E-05                  | 2.00E-03                 | 8.32E-03                         | 8.89E-04                          | 2.25E-02                         | N                    |
| Chlorobenzene               | 5.08E+02                 | ns       | 3.10E+02                 | n           | 5.38E-02                         | 6.80E-02                  | 7.50E-02                 | 6.06E-01                         | 7.65E-01                          | 8.44E-01                         | N                    |
| Chloroethane                | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Chloroform                  | 5.72E+00                 | c        | 3.00E-01                 | c           | 4.68E-04                         | 5.50E-05                  | -                        | 5.27E-03                         | 6.19E-04                          | -                                | N                    |
| Chloromethane               | 3.56E+01                 | c        | 1.20E+02                 | n           | 4.18E-03                         | 4.90E-02                  | -                        | 4.70E-02                         | 5.51E-01                          | -                                | N                    |
| 4-Chloro-3-methylphenol     | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 4-Chloroaniline             | -                        | -        | 2.40E+00                 | c           | -                                | 1.20E-04                  | -                        | -                                | 1.35E-03                          | -                                | N                    |
| 4-Chlorophenyl phenyl ether | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 4-Chlorotoluene             | -                        | -        | 5.50E+03                 | ns          | -                                | 2.80E+00                  | -                        | -                                | 3.15E+01                          | -                                | N                    |
| 2-Chloronaphthalene         | 6.26E+03                 | ns       | 6.30E+03                 | ns          | 1.35E+01                         | 1.80E+01                  | -                        | 1.52E+02                         | 2.03E+02                          | -                                | N                    |
| 2-Chlorophenol              | 3.91E+02                 | n        | 3.90E+02                 | n           | 1.53E-01                         | 2.00E-01                  | -                        | 1.72E+00                         | 2.25E+00                          | -                                | N                    |
| 2-Chlorotoluene             | 1.56E+03                 | ns       | 1.60E+03                 | ns          | 6.24E-01                         | 8.00E-01                  | -                        | 7.02E+00                         | 9.00E+00                          | -                                | N                    |
| Chromium III                | 1.13E+05                 | nl       | 1.20E+05                 | nm          | 9.86E+07                         | 9.90E+07                  | -                        | 1.11E+09                         | 1.11E+09                          | -                                | Y                    |
| Chromium VI                 | 2.19E+02                 | n        | 2.30E+02                 | n           | 2.11E+00                         | 2.10E+00                  | -                        | 2.37E+01                         | 2.36E+01                          | -                                | N                    |
| Chrysene                    | 4.81E+02                 | c        | 1.50E+01                 | c           | 3.26E+01                         | 1.40E+00                  | -                        | 3.67E+02                         | 1.58E+01                          | -                                | N                    |
| cis-1,2-DCE                 | 7.82E+02                 | n        | 7.80E+02                 | n           | 9.43E-02                         | 1.10E-01                  | 2.10E-02                 | 1.06E+00                         | 1.24E+00                          | 2.36E-01                         | N                    |
| cis-1,3-Dichloropropene     | 2.35E+01                 | c        | 1.70E+00                 | c*          | 1.35E-03                         | 1.60E-04                  | -                        | 1.52E-02                         | 1.80E-03                          | -                                | N                    |
| Cobalt                      | -                        | -        | 2.30E+01                 | n           | -                                | 4.90E-01                  | -                        | -                                | 5.51E+00                          | -                                | Y                    |
| Cyanide                     | 1.56E+03                 | n        | 1.60E+03                 | n           | 7.44E+00                         | 7.40E+00                  | 2.00E+00                 | 8.37E+01                         | 8.33E+01                          | 2.25E+01                         | N                    |
| 1,1-Dichloroethane          | 6.29E+01                 | c        | 3.40E+00                 | c           | 6.09E-03                         | 7.00E-04                  | -                        | 6.85E-02                         | 7.88E-03                          | -                                | N                    |
| 1,1-Dichloroethene          | 6.18E+02                 | n        | 2.50E+02                 | n           | 1.19E-01                         | 1.20E-01                  | 2.60E-03                 | 1.34E+00                         | 1.35E+00                          | 2.93E-02                         | N                    |
| 1,1-Dichloropropene         | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 1,2-Dibromo-3-chloropropane | 1.94E-01                 | c        | 5.60E-03                 | c           | 2.97E-06                         | 1.50E-07                  | 9.20E-05                 | 3.35E-05                         | 1.69E-06                          | 1.04E-03                         | N                    |
| 1,2-Dibromomethane (EDB)    | 5.74E-01                 | c        | 3.40E-02                 | c           | 1.58E-05                         | 1.90E-06                  | 1.50E-05                 | 1.78E-04                         | 2.14E-05                          | 1.69E-04                         | N                    |
| 1,2-Dichlorobenzene         | 3.01E+03                 | ns       | 2.00E+03                 | ns          | 3.13E-01                         | 4.00E-01                  | 6.60E-01                 | 3.53E+00                         | 4.50E+00                          | 7.43E+00                         | N                    |
| 3,3'-Dichlorobenzidine      | 8.71E+00                 | c        | 1.10E+00                 | c           | 1.70E-02                         | 2.30E-03                  | -                        | 1.92E-01                         | 2.59E-02                          | -                                | N                    |
| 1,2-Dichloroethane (EDC)    | 7.74E+00                 | c        | 4.50E-01                 | c           | 3.65E-04                         | 4.40E-05                  | 1.50E-03                 | 4.11E-03                         | 4.95E-04                          | 1.69E-02                         | N                    |
| 1,2-Dichloropropane         | 1.47E+01                 | c        | 9.30E-01                 | c*          | 1.11E-03                         | 1.30E-04                  | 1.70E-03                 | 1.25E-02                         | 1.46E-03                          | 1.91E-02                         | N                    |
| 1,3-Dichlorobenzene         | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 1,3-Dichloropropane         | -                        | -        | 1.60E+03                 | n           | -                                | 2.70E-01                  | -                        | -                                | 3.04E+00                          | -                                | N                    |
| 1,4-Dichlorobenzene         | 3.21E+01                 | c        | 2.60E+00                 | c           | 3.57E-03                         | 4.60E-04                  | 8.10E-02                 | 4.02E-02                         | 5.18E-03                          | 9.11E-01                         | N                    |
| 2,2-Dichloropropane         | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 2,4-Dichlorophenol          | 1.83E+02                 | n        | 1.80E+02                 | n           | 1.37E-01                         | 1.80E-01                  | -                        | 1.54E+00                         | 2.03E+00                          | -                                | N                    |
| 2,4-Dimethylphenol          | 1.22E+03                 | n        | 1.20E+03                 | n           | 9.12E-01                         | 1.20E+00                  | -                        | 1.03E+01                         | 1.35E+01                          | -                                | N                    |
| 4,6-Dinitro-2-methylphenol  | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 2,4-Dinitrophenol           | 1.22E+02                 | n        | 1.20E+02                 | n           | 5.25E-02                         | 6.80E-02                  | -                        | 5.91E-01                         | 7.65E-01                          | -                                | N                    |
| 2,4-Dinitrotoluene          | 1.26E+01                 | c        | 1.60E+00                 | c*          | 1.56E-03                         | 2.00E-04                  | -                        | 1.75E-02                         | 2.25E-03                          | -                                | N                    |
| 2,6-Dinitrotoluene          | 6.12E+01                 | n        | 6.10E+01                 | n           | 2.67E-02                         | 3.40E-02                  | -                        | 3.00E-01                         | 3.83E-01                          | -                                | N                    |
| Dibenz(a,h)anthracene       | 4.81E-01                 | c        | 1.50E-02                 | c           | 3.62E-01                         | 1.50E-02                  | -                        | 4.07E+00                         | 1.69E-01                          | -                                | N                    |
| Dibenzofuran                | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Dibromochloromethane        | 1.13E+01                 | c        | 7.00E-01                 | c           | 3.38E-04                         | 4.00E-05                  | -                        | 3.80E-03                         | 4.50E-04                          | -                                | N                    |
| Dibromomethane              | -                        | -        | 7.80E+02                 | n           | -                                | 9.10E-02                  | -                        | -                                | 1.02E+00                          | -                                | N                    |
| Dichlorodifluoromethane     | 4.81E+02                 | n        | 1.90E+02                 | n           | 7.23E-01                         | 6.10E-01                  | -                        | 8.14E+00                         | 6.86E+00                          | -                                | N                    |
| Diethyl phthalate           | 4.89E+04                 | n        | 4.90E+04                 | n           | 1.06E+01                         | 1.30E+01                  | -                        | 1.19E+02                         | 1.46E+02                          | -                                | N                    |
| Dimethyl phthalate          | 6.11E+05                 | nl       | -                        | -           | 8.36E+01                         | -                         | -                        | 9.40E+02                         | -                                 | -                                | N                    |
| Di-n-butyl phthalate        | 6.11E+03                 | n        | -                        | -           | 8.63E+00                         | -                         | -                        | 9.70E+01                         | -                                 | -                                | N                    |
| Di-n-octyl phthalate        | -                        | -        | -                        | -           | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Ethylbenzene                | 6.96E+01                 | c        | 5.70E+00                 | c           | 1.46E-02                         | 1.90E-03                  | 8.90E-01                 | 1.64E-01                         | 2.14E-02                          | 1.00E+01                         | N                    |
| Fluoranthene                | 2.29E+03                 | n        | 2.30E+03                 | n           | 1.55E+02                         | 2.10E+02                  | -                        | 1.75E+03                         | 2.36E+03                          | -                                | N                    |
| Fluorene                    | 2.29E+03                 | ns       | 2.30E+03                 | n           | 2.50E+01                         | 3.30E+01                  | -                        | 2.81E+02                         | 3.71E+02                          | -                                | N                    |
| Hexachlorobenzene           | 2.45E+00                 | c        | 3.00E-01                 | c           | 2.21E-03                         | 2.90E-04                  | 7.00E-03                 | 2.48E-02                         | 3.26E-03                          | 7.88E-02                         | N                    |

**TABLE 3**  
**Residential Soil Screening Levels**  
**Bloomfield Refinery - Bloomfield, New Mexico**

| Analyte                        | NMED                     |          | EPA                      |             | Cross Media Soil-to-Ground Water |          |                           | Cross Media Soil-to-Ground Water |          |          | Constituent Detected |
|--------------------------------|--------------------------|----------|--------------------------|-------------|----------------------------------|----------|---------------------------|----------------------------------|----------|----------|----------------------|
|                                | Residential Soil (mg/kg) | Endpoint | Residential Soil (mg/kg) | ResSoil key | NMED                             | EPA      | GW_Risk-based SSL (mg/kg) | GW_MCL-based SSL (mg/kg)         | NMED     | EPA      |                      |
| Applicable depth interval      | 0-10'                    |          | 0-2'                     |             | All depths                       |          |                           | All depths                       |          |          |                      |
| Hexachlorobutadiene            | -                        | -        | 6.20E+00                 | c**         | -                                | 1.90E-03 | -                         | -                                | 2.14E-02 | -        | N                    |
| Hexachlorocyclopentadiene      | 3.67E+02                 | n        | 3.70E+02                 | n           | 6.13E-01                         | 8.00E-01 | 1.80E-01                  | 6.90E+00                         | 9.00E+00 | 2.03E+00 | N                    |
| Hexachloroethane               | 6.11E+01                 | n        | 3.50E+01                 | c**         | 1.93E-02                         | 3.20E-03 | -                         | 2.17E-01                         | 3.60E-02 | -        | N                    |
| 2-Hexanone                     | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | Y                    |
| Indeno(1,2,3-cd)pyrene         | 4.81E+00                 | c        | 1.50E-01                 | c           | 3.70E+00                         | 1.60E-01 | -                         | 4.16E+01                         | 1.80E+00 | -        | N                    |
| Isophorone                     | 4.13E+03                 | c        | 5.10E+02                 | c*          | 1.85E-01                         | 2.20E-02 | -                         | 2.08E+00                         | 2.48E-01 | -        | N                    |
| Isopropylbenzene (cumene)      | 3.21E+03                 | ns       | 2.20E+03                 | ns          | 9.86E-01                         | 1.30E+00 | -                         | 1.11E+01                         | 1.46E+01 | -        | N                    |
| 4-Isopropyltoluene             | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | Y                    |
| Lead                           | 4.00E+02                 | IEUBK    | 4.00E+02                 | nL          | -                                | -        | -                         | -                                | -        | -        | Y                    |
| Mercury                        | 7.71E+00                 | ns       | 4.30E+00                 | ns          | 2.93E-02                         | 3.00E-02 | 1.00E-01                  | 3.30E-01                         | 3.38E-01 | 1.13E+00 | Y                    |
| Methyl tert-butyl ether (MTBE) | 8.62E+02                 | c        | 3.90E+01                 | c           | 2.29E-02                         | 2.70E-03 | -                         | 2.58E-01                         | 3.04E-02 | -        | N                    |
| Methylene chloride             | 1.99E+02                 | c        | 1.10E+01                 | c           | 1.07E-02                         | 1.20E-03 | 1.30E-03                  | 1.21E-01                         | 1.35E-02 | 1.46E-02 | Y                    |
| 1-Methylnaphthalene            | -                        | -        | 2.20E+01                 | c           | -                                | 1.50E-02 | -                         | -                                | 1.69E-01 | -        | N                    |
| 2-Methylnaphthalene            | -                        | -        | 3.10E+02                 | n           | -                                | 9.00E-01 | -                         | -                                | 1.01E+01 | -        | N                    |
| 2-Methylphenol                 | -                        | -        | 3.10E+03                 | n           | -                                | 2.00E+00 | -                         | -                                | 2.25E+01 | -        | N                    |
| 3+4-Methylphenol               | -                        | -        | 3.10E+02                 | n           | -                                | 1.90E-01 | -                         | -                                | 2.14E+00 | -        | Y                    |
| 4-Methyl-2-pentanone           | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | Y                    |
| 2-Nitroaniline                 | -                        | -        | 1.80E+02                 | n           | -                                | 3.30E-02 | -                         | -                                | 3.71E-01 | -        | N                    |
| 3-Nitroaniline                 | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | N                    |
| 4-Nitroaniline                 | -                        | -        | 2.40E+01                 | c*          | -                                | 1.00E-03 | -                         | -                                | 1.13E-02 | -        | N                    |
| 2-Nitrophenol                  | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | N                    |
| 4-Nitrophenol                  | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | N                    |
| Naphthalene                    | 4.50E+01                 | c        | 3.90E+00                 | c*          | 4.19E-03                         | 5.50E-04 | -                         | 4.72E-02                         | 6.19E-03 | -        | N                    |
| n-Butylbenzene                 | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | N                    |
| Nickel                         | 1.56E+03                 | n        | 1.40E+04                 | c           | 4.77E+01                         | 4.80E+01 | -                         | 5.36E+02                         | 5.40E+02 | -        | Y                    |
| Nitrobenzene                   | 4.94E+01                 | c        | 4.40E+00                 | c*          | 6.86E-03                         | 7.10E-05 | -                         | 7.72E-02                         | 7.99E-04 | -        | N                    |
| N-Nitrosodi-n-propylamine      | -                        | -        | 6.90E-02                 | c           | -                                | 1.10E-05 | -                         | -                                | 1.24E-04 | -        | N                    |
| N-Nitrosodiphenylamine         | 8.00E+02                 | c        | 9.90E+01                 | c           | 1.29E+00                         | 1.70E-01 | -                         | 1.45E+01                         | 1.91E+00 | -        | N                    |
| n-Propylbenzene                | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | N                    |
| Pentachlorophenol              | 2.07E+01                 | c        | 3.00E+00                 | c           | 2.94E-02                         | 3.90E-03 | 7.00E-03                  | 3.30E-01                         | 4.39E-02 | 7.88E-02 | N                    |
| Phenanthrene                   | 1.83E+03                 | ns       | -                        | -           | 8.34E+01                         | -        | -                         | 9.39E+02                         | -        | -        | N                    |
| Phenol                         | 1.83E+04                 | n        | 1.80E+04                 | n           | 6.30E+00                         | 8.10E+00 | -                         | 7.09E+01                         | 9.11E+01 | -        | N                    |
| Pyrene                         | 1.72E+03                 | ns       | 1.70E+03                 | n           | 1.12E+02                         | 1.50E+02 | -                         | 1.26E+03                         | 1.69E+03 | -        | N                    |
| Pyridine                       | -                        | -        | 7.80E+01                 | n           | -                                | 9.70E-03 | -                         | -                                | 1.09E-01 | -        | N                    |
| sec-Butylbenzene               | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | N                    |
| Selenium                       | 3.91E+02                 | n        | 3.90E+02                 | n           | 9.65E-01                         | 9.50E-01 | 2.60E-01                  | 1.09E+01                         | 1.07E+01 | 2.93E+00 | N                    |
| Silver                         | 3.91E+02                 | n        | 3.90E+02                 | n           | 1.57E+00                         | 1.60E+00 | -                         | 1.76E+01                         | 1.80E+01 | -        | Y                    |
| Styrene                        | 8.97E+03                 | ns       | 6.50E+03                 | ns          | 1.56E+00                         | 2.00E+00 | 1.20E-01                  | 1.76E+01                         | 2.25E+01 | 1.35E+00 | N                    |
| 1,2,3-Trichlorobenzene         | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | N                    |
| 1,1,1,2-Tetrachloroethane      | 2.92E+01                 | c        | 2.00E+00                 | c           | 1.73E-03                         | 2.10E-04 | -                         | 1.94E-02                         | 2.36E-03 | -        | N                    |
| 1,1,1-Trichloroethane          | 2.18E+04                 | ns       | 9.00E+03                 | ns          | 2.98E+00                         | 3.30E+00 | 7.20E-02                  | 3.35E+01                         | 3.71E+01 | 8.10E-01 | N                    |
| 1,1,2,2-Tetrachloroethane      | 7.97E+00                 | c        | 5.90E-01                 | c           | 2.25E-04                         | 2.80E-05 | -                         | 2.53E-03                         | 3.15E-04 | -        | N                    |
| 1,1,2-Trichloroethane          | 1.72E+01                 | c        | 1.10E+00                 | c           | 6.74E-04                         | 8.20E-05 | 1.70E-03                  | 7.58E-03                         | 9.23E-04 | 1.91E-02 | N                    |
| 2,4,5-Trichlorophenol          | 6.11E+03                 | n        | 6.10E+03                 | n           | 7.13E+00                         | 9.40E+00 | -                         | 8.02E+01                         | 1.06E+02 | -        | N                    |
| 2,4,6-Trichlorophenol          | 6.11E+01                 | n        | 4.40E+01                 | c**         | 7.13E-02                         | 1.60E-02 | -                         | 8.02E-01                         | 1.80E-01 | -        | N                    |
| 1,2,3-Trichloropropane         | 9.15E-01                 | c        | 9.10E-02                 | c           | 3.56E-05                         | 4.40E-06 | -                         | 4.01E-04                         | 4.95E-05 | -        | N                    |
| 1,2,4-Trichlorobenzene         | 1.43E+02                 | ns       | 8.70E+01                 | n           | 1.02E-02                         | 1.30E-02 | 1.10E-01                  | 1.15E-01                         | 1.46E-01 | 1.24E+00 | N                    |
| 1,2,4-Trimethylbenzene         | -                        | -        | 6.70E+01                 | n           | -                                | 2.40E-02 | -                         | -                                | 2.70E-01 | -        | N                    |
| 1,3,5-Trimethylbenzene         | -                        | -        | 4.70E+01                 | n           | -                                | 2.00E-02 | -                         | -                                | 2.25E-01 | -        | N                    |
| tert-Butylbenzene              | -                        | -        | -                        | -           | -                                | -        | -                         | -                                | -        | -        | N                    |
| Tetrachloroethene (PCE)        | 6.99E+00                 | c        | 5.70E-01                 | c           | 4.49E-04                         | 5.20E-05 | 2.40E-03                  | 5.05E-03                         | 5.85E-04 | 2.70E-02 | N                    |
| Toluene                        | 5.57E+03                 | ns       | 5.00E+03                 | ns          | 1.38E+00                         | 1.70E+00 | 7.60E-01                  | 1.56E+01                         | 1.91E+01 | 8.55E+00 | Y                    |
| trans-1,2-DCE                  | 2.73E+02                 | n        | 1.10E+02                 | n           | 3.01E-02                         | 3.40E-02 | 3.20E-02                  | 3.39E-01                         | 3.83E-01 | 3.60E-01 | N                    |
| trans-1,3-Dichloropropene      | 2.35E+01                 | c        | 1.70E+00                 | c*          | 1.35E-03                         | 1.60E-04 | -                         | 1.52E-02                         | 1.80E-03 | -        | N                    |
| Trichloroethene (TCE)          | 4.57E+01                 | c        | 2.80E+00                 | c           | 5.30E-03                         | 6.10E-04 | 1.90E-03                  | 5.96E-02                         | 6.86E-03 | 2.14E-02 | N                    |
| Trichlorofluoromethane         | 2.01E+03                 | ns       | 8.00E+02                 | n           | 9.01E-01                         | 8.40E-01 | -                         | 1.01E+01                         | 9.45E+00 | -        | N                    |
| Vanadium                       | 3.91E+02                 | n        | 5.50E+02                 | n           | 1.83E+02                         | 2.60E+02 | -                         | 2.05E+03                         | 2.93E+03 | -        | Y                    |
| Vinyl chloride                 | 8.65E-01                 | c        | 6.00E-02                 | c           | 2.88E-04                         | 5.60E-06 | 7.00E-04                  | 3.24E-03                         | 6.30E-05 | 7.88E-03 | N                    |
| Xylenes, Total                 | 1.09E+03                 | ns       | 6.00E+02                 | ns          | 1.76E-01                         | 2.30E-01 | 1.10E+01                  | 1.98E+00                         | 2.59E+00 | 1.24E+02 | N                    |
| Zinc                           | 2.35E+04                 | n        | 2.30E+04                 | n           | 6.82E+02                         | 6.80E+02 | -                         | 7.67E+03                         | 7.65E+03 | -        | Y                    |

c - carcinogen  
n - noncarcinogen  
cs - carcinogen, SSL may exceed saturation  
ns - noncarcinogen, SSL may exceed saturation  
nl - noncarcinogen, SSL may exceed ceiling limit  
nls - noncarcinogen, SSL may exceed both saturation and ceiling limit  
- no screening value currently available

NMED - Technical Background Document for Development of Soil Screening Levels - Revision 5.0 ( August 2009)  
EPA - Regional Screening Levels (April 2009)

**TABLE 4**  
**Non- Residential Soil Screening Levels**  
**Bloomfield Refinery - Bloomfield, New Mexico**

| Analyte                     | NMED               |                       |                  |                          | EPA                |              | Cross Media Soil-to-Ground-Water |                           |                          | Cross Media Soil-to-Ground-Water |                                   |                                  | Constituent Detected |
|-----------------------------|--------------------|-----------------------|------------------|--------------------------|--------------------|--------------|----------------------------------|---------------------------|--------------------------|----------------------------------|-----------------------------------|----------------------------------|----------------------|
|                             | IndOccSoil (mg/kg) | IndOccSoil (Endpoint) | ConsWork (mg/kg) | ConsWork Soil (Endpoint) | Industrial (mg/kg) | IndSoil _key | NMED                             | EPA                       |                          | NMED                             | EPA                               |                                  |                      |
|                             |                    |                       |                  |                          |                    |              | DAF1 (mg/kg)                     | GW_Risk-based SSL (mg/kg) | GW_MCL-based SSL (mg/kg) | DAF (11.25) (mg/kg)              | GW_Risk-based SSL (11.25) (mg/kg) | GW_MCL-based SSL (11.25) (mg/kg) |                      |
| Applicable depth interval   | 0-2'               |                       | 0-10'            |                          | 0-2'               |              | All depths                       |                           |                          | All depths                       |                                   |                                  |                      |
| Acenaphthene                | 3.67E+04           | ns                    | 1.86E+04         | n                        | 3.30E+04           | n            | 2.05E+01                         | 2.70E+01                  | -                        | 2.31E+02                         | 3.04E+02                          | -                                | N                    |
| Acenaphthylene              | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Acetone                     | 8.51E+05           | nls                   | 2.63E+05         | nls                      | 6.10E+05           | nms          | 3.84E+00                         | 4.40E+00                  | -                        | 4.32E+01                         | 4.95E+01                          | -                                | Y                    |
| Aniline                     | -                  | -                     | -                | -                        | 3.00E+02           | c*           | -                                | 3.40E-03                  | -                        | -                                | 3.83E-02                          | -                                | N                    |
| Anthracene                  | 1.83E+05           | nl                    | 6.68E+04         | ns                       | 1.70E+05           | nm           | 3.37E+02                         | 4.50E+02                  | -                        | 3.79E+03                         | 5.06E+03                          | -                                | N                    |
| Antimony                    | 4.54E+02           | n                     | 1.24E+02         | n                        | 4.10E+02           | n            | 6.61E-01                         | 6.60E-01                  | 2.70E-01                 | 7.44E+00                         | 7.43E+00                          | 3.04E+00                         | N                    |
| Arsenic                     | 1.77E+01           | c                     | 6.54E+01         | n                        | 1.60E+00           | c            | 1.31E-02                         | 1.30E-03                  | 2.90E-01                 | 1.48E-01                         | 1.46E-02                          | 3.26E+00                         | Y                    |
| Azobenzene                  | -                  | -                     | -                | -                        | 2.20E+01           | c            | -                                | 5.10E-04                  | -                        | -                                | 5.74E-03                          | -                                | N                    |
| Barium                      | 2.24E+05           | nl                    | 4.35E+03         | n                        | 1.90E+05           | nm           | 3.01E+02                         | 3.00E+02                  | 8.20E+01                 | 3.39E+03                         | 3.38E+03                          | 9.23E+02                         | Y                    |
| Beryllium                   | 2.26E+03           | n                     | 1.44E+02         | n                        | 2.00E+03           | n            | 5.77E+01                         | 5.80E+01                  | 3.20E+00                 | 6.49E+02                         | 6.53E+02                          | 3.60E+01                         | Y                    |
| Benzo(a)anthracene          | 2.34E+01           | c                     | 2.13E+02         | c                        | 2.10E+00           | c            | 3.20E-01                         | 1.40E-02                  | -                        | 3.59E+00                         | 1.58E-01                          | -                                | N                    |
| Benzene                     | 8.54E+01           | c                     | 4.71E+02         | n                        | 5.60E+00           | c*           | 1.85E-03                         | 2.30E-04                  | 2.80E-03                 | 2.08E-02                         | 2.59E-03                          | 3.15E-02                         | N                    |
| Benzo(a)pyrene              | 2.34E+00           | c                     | 2.13E+01         | c                        | 2.10E+01           | c            | 1.09E-01                         | 4.60E-03                  | 3.10E-01                 | 1.22E+00                         | 5.18E-02                          | 3.49E+00                         | N                    |
| Benzo(b)fluoranthene        | 2.34E+01           | c                     | 2.13E+02         | c                        | 2.10E+00           | c            | 1.11E+00                         | 4.70E-02                  | -                        | 1.25E+01                         | 5.29E-01                          | -                                | N                    |
| Benzo(g,h,i)perylene        | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Benzo(k)fluoranthene        | 2.34E+02           | c                     | 2.06E+03         | c                        | 2.10E+01           | c            | 1.09E+01                         | 4.60E-01                  | -                        | 1.22E+02                         | 5.18E+00                          | -                                | Y                    |
| Benzoic acid                | -                  | -                     | -                | -                        | 2.50E+06           | nm           | -                                | 3.30E+01                  | -                        | -                                | 3.71E+02                          | -                                | N                    |
| Benzyl alcohol              | -                  | -                     | -                | -                        | 3.10E+05           | nm           | -                                | 4.20E+00                  | -                        | -                                | 4.73E+01                          | -                                | Y                    |
| Bis(2-chloroethoxy)methane  | -                  | -                     | -                | -                        | 1.80E+03           | n            | -                                | 2.30E-02                  | -                        | -                                | 2.59E-01                          | -                                | N                    |
| Bis(2-chloroethyl)ether     | 1.36E+01           | c                     | 1.47E+02         | c                        | 9.00E-01           | c            | 2.33E-05                         | 2.70E-06                  | -                        | 2.62E-04                         | 3.04E-05                          | -                                | N                    |
| Bis(2-chloroisopropyl)ether | 4.54E+02           | c                     | 3.10E+03         | cs                       | -                  | -            | 2.56E-03                         | -                         | -                        | 2.88E-02                         | -                                 | -                                | N                    |
| Bis(2-ethylhexyl)phthalate  | 1.37E+03           | c                     | 4.76E+03         | n                        | 1.20E+02           | c            | 1.19E+01                         | 1.60E+00                  | 2.00E+00                 | 1.34E+02                         | 1.80E+01                          | 2.25E+01                         | N                    |
| Bromobenzene                | -                  | -                     | -                | -                        | 4.10E+02           | n            | -                                | 1.50E-02                  | -                        | -                                | 1.69E-01                          | -                                | N                    |
| Bromodichloromethane        | 2.92E+01           | c                     | 3.50E+03         | cs                       | 1.40E+00           | c            | 2.76E-04                         | 3.30E-05                  | -                        | 3.11E-03                         | 3.71E-04                          | -                                | N                    |
| Bromoform                   | -                  | -                     | -                | -                        | 2.20E+02           | c*           | -                                | 2.30E-03                  | -                        | -                                | 2.59E-02                          | -                                | N                    |
| Bromomethane                | 8.36E+01           | n                     | 6.71E+01         | n                        | 3.50E+01           | n            | 1.94E-03                         | 2.20E-03                  | -                        | 2.18E-02                         | 2.48E-02                          | -                                | N                    |
| 4-Bromophenyl phenyl ether  | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 2-Butanone (MEK)            | 3.69E+05           | nl                    | 1.48E+05         | nls                      | 1.90E+05           | nms          | 1.27E+00                         | 1.50E+00                  | -                        | 1.43E+01                         | 1.69E+01                          | -                                | Y                    |
| Butyl benzyl phthalate      | -                  | -                     | -                | -                        | 9.10E+02           | c            | -                                | 6.70E-01                  | -                        | -                                | 7.54E+00                          | -                                | N                    |
| Cadmium                     | 1.12E+03           | n                     | 3.09E+02         | n                        | 8.00E+02           | n            | 1.37E+00                         | 1.40E+00                  | 3.80E-01                 | 1.55E+01                         | 1.58E+01                          | 4.28E+00                         | N                    |
| Carbazole                   | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Carbon disulfide            | 7.54E+03           | ns                    | 5.89E+03         | ns                       | 3.00E+03           | ns           | 2.52E-01                         | 2.70E-01                  | -                        | 2.84E+00                         | 3.04E+00                          | -                                | Y                    |
| Carbon tetrachloride        | 2.43E+01           | c                     | 1.99E+02         | n                        | 1.30E+00           | c            | 7.39E-04                         | 7.90E-05                  | 2.00E-03                 | 8.32E-03                         | 8.89E-04                          | 2.25E-02                         | N                    |
| Chlorobenzene               | 2.14E+03           | n                     | 1.58E+03         | ns                       | 1.50E+03           | ns           | 5.38E-02                         | 6.80E-02                  | 7.50E-02                 | 6.06E-01                         | 7.65E-01                          | 8.44E-01                         | N                    |
| Chloroethane                | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Chloroform                  | 3.19E+01           | c                     | 6.71E+02         | c                        | 1.50E+00           | c            | 4.68E-04                         | 5.50E-05                  | -                        | 5.27E-03                         | 6.19E-04                          | -                                | N                    |
| Chloromethane               | 1.98E+02           | c                     | 1.13E+03         | n                        | 5.10E+02           | n            | 4.18E-03                         | 4.90E-02                  | -                        | 4.70E-02                         | 5.51E-01                          | -                                | N                    |
| 4-Chloro-3-methylphenol     | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 4-Chloroaniline             | -                  | -                     | -                | -                        | 8.60E+00           | c            | -                                | 1.20E-04                  | -                        | -                                | 1.35E-03                          | -                                | N                    |
| 4-Chlorophenyl phenyl ether | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 4-Chlorotoluene             | -                  | -                     | -                | -                        | 7.20E+04           | ns           | -                                | 2.80E+00                  | -                        | -                                | 3.15E+01                          | -                                | N                    |
| 2-Chloronaphthalene         | 9.08E+04           | ns                    | 2.48E+04         | ns                       | 8.20E+04           | ns           | 1.35E+01                         | 1.80E+01                  | -                        | 1.52E+02                         | 2.03E+02                          | -                                | N                    |
| 2-Chlorophenol              | 5.68E+03           | n                     | 1.55E+03         | n                        | 5.10E+03           | n            | 1.53E-01                         | 2.00E-01                  | -                        | 1.72E+00                         | 2.25E+00                          | -                                | N                    |
| 2-Chlorotoluene             | 2.27E+04           | ns                    | 6.19E+03         | ns                       | 2.00E+04           | ns           | 6.24E-01                         | 8.00E-01                  | -                        | 7.02E+00                         | 9.00E+00                          | -                                | N                    |
| Chromium III                | 1.57E+06           | nl                    | 4.47E+05         | nl                       | 1.50E+06           | nm           | 9.86E+07                         | 9.90E+07                  | -                        | 1.11E+09                         | 1.11E+09                          | -                                | Y                    |
| Chromium VI                 | 2.92E+03           | n                     | 4.49E+02         | n                        | 1.40E+03           | c            | 2.11E+00                         | 2.10E+00                  | -                        | 2.37E+01                         | 2.36E+01                          | -                                | N                    |
| Chrysene                    | 2.34E+03           | c                     | 2.06E+04         | c                        | 2.10E+02           | c            | 3.26E+01                         | 1.40E+00                  | -                        | 3.67E+02                         | 1.58E+01                          | -                                | N                    |
| cis-1,2-DCE                 | 1.14E+04           | ns                    | 3.10E+03         | cs                       | 1.00E+04           | ns           | 9.43E-02                         | 1.10E-01                  | 2.10E-02                 | 1.06E+00                         | 1.24E+00                          | 2.36E-01                         | N                    |
| cis-1,3-Dichloropropene     | 1.26E+02           | c                     | 5.10E+02         | n                        | 8.40E+00           | c*           | 1.35E-03                         | 1.60E-04                  | -                        | 1.52E-02                         | 1.80E-03                          | -                                | N                    |
| Cobalt                      | -                  | -                     | -                | -                        | 3.00E+02           | n            | -                                | 4.90E-01                  | -                        | -                                | 5.51E+00                          | -                                | Y                    |
| Cyanide                     | 2.27E+04           | n                     | 6.19E+03         | n                        | 2.00E+04           | n            | 7.44E+00                         | 7.40E+00                  | 2.00E+00                 | 8.37E+01                         | 8.33E+01                          | 2.25E+01                         | N                    |
| 1,1-Dichloroethane          | 3.50E+02           | c                     | 6.88E+03         | cs                       | 1.70E+01           | c            | 6.09E-03                         | 7.00E-04                  | -                        | 6.85E-02                         | 7.88E-03                          | -                                | N                    |
| 1,1-Dichloroethene          | 2.22E+03           | ns                    | 1.83E+03         | ns                       | 1.10E+03           | n            | 1.19E-01                         | 1.20E-01                  | 2.60E-03                 | 1.34E+00                         | 1.35E+00                          | 2.93E-02                         | N                    |
| 1,1-Dichloropropene         | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 1,2-Dibromo-3-chloropropane | 1.09E+00           | c                     | 2.30E+01         | c                        | 7.30E-02           | c            | 2.97E-06                         | 1.50E-07                  | 9.20E-05                 | 3.35E-05                         | 1.69E-06                          | 1.04E-03                         | N                    |
| 1,2-Dibromoethane (EDB)     | 3.14E+00           | c                     | 4.86E+01         | c                        | 1.70E-01           | c            | 1.58E-05                         | 1.90E-06                  | 1.50E-05                 | 1.78E-04                         | 2.14E-05                          | 1.69E-04                         | N                    |
| 1,2-Dichlorobenzene         | 1.43E+04           | ns                    | 9.71E+03         | ns                       | 1.00E+04           | ns           | 3.13E-01                         | 4.00E-01                  | 6.60E-01                 | 3.53E+00                         | 4.50E+00                          | 7.43E+00                         | N                    |
| 3,3'-Dichlorobenzidine      | 4.26E+01           | c                     | 3.71E+02         | c                        | 3.80E+00           | c            | 1.70E-02                         | 2.30E-03                  | -                        | 1.92E-01                         | 2.59E-02                          | -                                | N                    |
| 1,2-Dichloroethane (EDC)    | 4.28E+01           | c                     | 7.51E+02         | c                        | 2.20E+00           | c            | 3.65E-04                         | 4.40E-05                  | 1.50E-03                 | 4.11E-03                         | 4.95E-04                          | 1.69E-02                         | N                    |
| 1,2-Dichloropropane         | 8.17E+01           | c                     | 1.17E+02         | n                        | 4.70E+00           | c*           | 1.11E-03                         | 1.30E-04                  | 1.70E-03                 | 1.25E-02                         | 1.46E-03                          | 1.91E-02                         | N                    |
| 1,3-Dichlorobenzene         | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 1,3-Dichloropropane         | -                  | -                     | -                | -                        | 2.00E+04           | ns           | -                                | 2.70E-01                  | -                        | -                                | 3.04E+00                          | -                                | N                    |
| 1,4-Dichlorobenzene         | 1.80E+02           | c                     | 3.78E+03         | cs                       | 1.30E+01           | c            | 3.57E-03                         | 4.60E-04                  | 8.10E-02                 | 4.02E-02                         | 5.18E-03                          | 9.11E-01                         | N                    |
| 2,2-Dichloropropane         | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 2,4-Dichlorophenol          | 2.05E+03           | n                     | 7.15E+02         | n                        | 1.80E+03           | n            | 1.37E-01                         | 1.80E-01                  | -                        | 1.54E+00                         | 2.03E+00                          | -                                | N                    |
| 2,4-Dimethylphenol          | 1.37E+04           | n                     | 4.76E+03         | n                        | 1.20E+04           | n            | 9.12E-01                         | 1.20E+00                  | -                        | 1.03E+01                         | 1.35E+01                          | -                                | N                    |
| 4,6-Dinitro-2-methylphenol  | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 2,4-Dinitrophenol           | 1.37E+03           | n                     | 4.76E+02         | n                        | 1.20E+03           | n            | 5.25E-02                         | 6.80E-02                  | -                        | 5.91E-01                         | 7.65E-01                          | -                                | N                    |
| 2,4-Dinitrotoluene          | 1.03E+02           | c                     | 4.76E+02         | n                        | 5.50E+00           | c            | 1.56E-03                         | 2.00E-04                  | -                        | 1.75E-02                         | 2.25E-03                          | -                                | N                    |
| 2,6-Dinitrotoluene          | 6.87E+02           | n                     | 2.39E+02         | n                        | 6.20E+02           | n            | 2.67E-02                         | 3.40E-02                  | -                        | 3.00E-01                         | 3.83E-01                          | -                                | N                    |
| Dibenz(a,h)anthracene       | 2.34E+00           | c                     | 2.13E+01         | c                        | 2.10E-01           | c            | 3.62E-01                         | 1.50E-02                  | -                        | 4.07E+00                         | 1.69E-01                          | -                                | N                    |
| Dibenzofuran                | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Dibromochloromethane        | 6.13E+01           | c                     | 1.99E+03         | c                        | 3.40E+00           | c            | 3.38E-04                         | 4.00E-05                  | -                        | 3.80E-03                         | 4.50E-04                          | -                                | N                    |
| Dibromomethane              | -                  | -                     | -                | -                        | 1.00E+04           | ns           | -                                | 9.10E-02                  | -                        | -                                | 1.02E+00                          | -                                | N                    |
| Dichlorodifluoromethane     | 1.55E+03           | ns                    | 1.37E+03         | ns                       | 7.80E+02           | n            | 7.23E-01                         | 6.10E-01                  | -                        | 8.14E+00                         | 6.86E+00                          | -                                | N                    |
| Diethyl phthalate           | 5.47E+05           | nl                    | 1.91E+05         | nl                       | 4.90E+05           | nm           | 1.06E+01                         | 1.30E+01                  | -                        | 1.19E+02                         | 1.46E+02                          | -                                | N                    |
| Dimethyl phthalate          | 6.84E+06           | nl                    | 2.38E+06         | nl                       | -                  | -            | 8.36E+01                         | -                         | -                        | 9.40E+02                         | -                                 | -                                | N                    |
| Di-n-butyl phthalate        | 6.84E+04           | n                     | 2.38E+04         | n                        | -                  | -            | 8.63E+00                         | -                         | -                        | 9.70E+01                         | -                                 | -                                | N                    |



**TABLE 4**  
**Non- Residential Soil Screening Levels**  
**Bloomfield Refinery - Bloomfield, New Mexico**

| Analyte                        | NMED               |                       |                  |                          | EPA                |              | Cross Media Soil-to-Ground Water |                           |                          | Cross Media Soil-to-Ground Water |                                   |                                  | Constituent Detected |
|--------------------------------|--------------------|-----------------------|------------------|--------------------------|--------------------|--------------|----------------------------------|---------------------------|--------------------------|----------------------------------|-----------------------------------|----------------------------------|----------------------|
|                                | IndOccSoil (mg/kg) | IndOccSoil (Endpoint) | ConsWork (mg/kg) | ConsWork Soil (Endpoint) | Industrial (mg/kg) | IndSoil _key | DAF1 (mg/kg)                     | GW_Risk-based SSL (mg/kg) | GW_MCL-based SSL (mg/kg) | DAF (11.25) (mg/kg)              | GW_Risk-based SSL (11.25) (mg/kg) | GW_MCL-based SSL (11.25) (mg/kg) |                      |
|                                | 0-2'               | 0-2'                  | 0-10'            | 0-10'                    | 0-2'               |              | All depths                       | All depths                | All depths               | All depths                       | All depths                        | All depths                       |                      |
| Di-n-octyl phthalate           | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Ethylbenzene                   | 3.85E+02           | c                     | 6.63E+03         | cs                       | 2.90E+01           | c            | 1.46E-02                         | 1.90E-03                  | 8.90E-01                 | 1.64E-01                         | 2.14E-02                          | 1.00E+01                         | N                    |
| Fluoranthene                   | 2.44E+04           | n                     | 8.91E+03         | n                        | 2.20E+04           | n            | 1.55E+02                         | 2.10E+02                  | -                        | 1.75E+03                         | 2.36E+03                          | -                                | N                    |
| Fluorene                       | 2.44E+04           | ns                    | 8.91E+03         | ns                       | 2.20E+04           | n            | 2.50E+01                         | 3.30E+01                  | -                        | 2.81E+02                         | 3.71E+02                          | -                                | N                    |
| Hexachlorobenzene              | 1.20E+01           | c                     | 1.03E+02         | c                        | 1.10E+00           | c            | 2.21E-03                         | 2.90E-04                  | 7.00E-03                 | 2.48E-02                         | 3.26E-03                          | 7.88E-02                         | N                    |
| Hexachlorobutadiene            | -                  | -                     | -                | -                        | 2.20E+01           | c*           | -                                | 1.90E-03                  | -                        | -                                | 2.14E-02                          | -                                | N                    |
| Hexachlorocyclopentadiene      | 4.10E+03           | n                     | 8.11E+02         | n                        | 3.70E+03           | n            | 6.13E-01                         | 8.00E-01                  | 1.80E-01                 | 6.90E+00                         | 9.00E+00                          | 2.03E+00                         | N                    |
| Hexachloroethane               | 6.84E+02           | n                     | 2.38E+02         | n                        | 1.20E+02           | c**          | 1.93E-02                         | 3.20E-03                  | -                        | 2.17E-01                         | 3.60E-02                          | -                                | N                    |
| 2-Hexanone                     | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | Y                    |
| Indeno(1,2,3-cd)pyrene         | 2.34E+01           | c                     | 2.13E+02         | c                        | 2.10E+00           | c            | 3.70E+00                         | 1.60E-01                  | -                        | 4.16E+01                         | 1.80E+00                          | -                                | N                    |
| Isophorone                     | 2.02E+04           | c                     | 4.75E+04         | n                        | 1.80E+03           | c*           | 1.85E-01                         | 2.20E-02                  | -                        | 2.08E+00                         | 2.48E-01                          | -                                | N                    |
| Isopropylbenzene (cumene)      | 1.49E+04           | ns                    | 1.03E+04         | ns                       | 1.10E+04           | ns           | 9.86E-01                         | 1.30E+00                  | -                        | 1.11E+01                         | 1.46E+01                          | -                                | N                    |
| 4-Isopropyltoluene             | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | Y                    |
| Lead                           | 8.00E+02           | IEUBK                 | 8.00E+02         | IEUBK                    | 8.00E+02           | nL           | -                                | -                         | -                        | -                                | -                                 | -                                | Y                    |
| Mercury                        | 4.99E+01           | n                     | 6.36E+01         | ns                       | 2.40E+01           | ns           | 2.93E-02                         | 3.00E-02                  | 1.00E-01                 | 3.30E-01                         | 3.38E-01                          | 1.13E+00                         | Y                    |
| Methyl tert-butyl ether (MTBE) | 4.69E+03           | c                     | 6.55E+04         | cs                       | 1.90E+02           | c            | 2.29E-02                         | 2.70E-03                  | -                        | 2.58E-01                         | 3.04E-02                          | -                                | N                    |
| Methylene chloride             | 1.09E+03           | c                     | 1.06E+04         | ns                       | 5.40E+01           | c            | 1.07E-02                         | 1.20E-03                  | 1.30E-03                 | 1.21E-01                         | 1.35E-02                          | 1.48E-02                         | Y                    |
| 1-Methylnaphthalene            | -                  | -                     | -                | -                        | 9.90E+01           | c            | -                                | 1.50E-02                  | -                        | -                                | 1.69E-01                          | -                                | N                    |
| 2-Methylnaphthalene            | -                  | -                     | -                | -                        | 4.10E+03           | ns           | -                                | 9.00E-01                  | -                        | -                                | 1.01E+01                          | -                                | N                    |
| 2-Methylphenol                 | -                  | -                     | -                | -                        | 3.10E+04           | n            | -                                | 2.00E+00                  | -                        | -                                | 2.25E+01                          | -                                | N                    |
| 3+4-Methylphenol               | -                  | -                     | -                | -                        | 3.10E+03           | n            | -                                | 1.90E-01                  | -                        | -                                | 2.14E+00                          | -                                | Y                    |
| 4-Methyl-2-pentanone           | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | Y                    |
| 2-Nitroaniline                 | -                  | -                     | -                | -                        | 1.80E+03           | n            | -                                | 3.30E-02                  | -                        | -                                | 3.71E-01                          | -                                | N                    |
| 3-Nitroaniline                 | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 4-Nitroaniline                 | -                  | -                     | -                | -                        | 8.60E+01           | c*           | -                                | 1.00E-03                  | -                        | -                                | 1.13E-02                          | -                                | N                    |
| 2-Nitrophenol                  | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 4-Nitrophenol                  | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Naphthalene                    | 2.52E+02           | cs                    | 7.02E+02         | ns                       | 2.00E+01           | c*           | 4.19E-03                         | 5.50E-04                  | -                        | 4.72E-02                         | 6.19E-03                          | -                                | N                    |
| n-Butylbenzene                 | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Nickel                         | 2.27E+04           | n                     | 6.19E+03         | n                        | 6.90E+04           | c            | 4.77E+01                         | 4.80E+01                  | -                        | 5.36E+02                         | 5.40E+02                          | -                                | Y                    |
| Nitrobenzene                   | 2.77E+02           | c                     | 5.20E+02         | n                        | 2.20E+01           | c*           | 6.86E-03                         | 7.10E-05                  | -                        | 7.72E-02                         | 7.99E-04                          | -                                | N                    |
| N-Nitrosodi-n-propylamine      | -                  | -                     | -                | -                        | 2.50E-01           | c            | -                                | 1.10E-05                  | -                        | -                                | 1.24E-04                          | -                                | N                    |
| N-Nitrosodiphenylamine         | 3.91E+03           | c                     | 3.40E+04         | c                        | 3.50E+02           | c            | 1.29E+00                         | 1.70E-01                  | -                        | 1.45E+01                         | 1.91E+00                          | -                                | N                    |
| n-Propylbenzene                | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Pentachlorophenol              | 1.00E+02           | c                     | 1.03E+03         | c                        | 9.00E+00           | c            | 2.94E-02                         | 3.90E-03                  | 7.00E-03                 | 3.30E-01                         | 4.39E-02                          | 7.88E-02                         | N                    |
| Phenanthrene                   | 2.05E+04           | ns                    | 7.15E+03         | ns                       | -                  | -            | 8.34E+01                         | -                         | -                        | 9.39E+02                         | -                                 | -                                | N                    |
| Phenol                         | 2.05E+05           | nl                    | 6.88E+04         | n                        | 1.80E+05           | nm           | 6.30E+00                         | 8.10E+00                  | -                        | 7.09E+01                         | 9.11E+01                          | -                                | N                    |
| Pyrene                         | 1.83E+04           | ns                    | 6.68E+03         | ns                       | 1.70E+04           | n            | 1.12E+02                         | 1.50E+02                  | -                        | 1.26E+03                         | 1.69E+03                          | -                                | N                    |
| Pyridine                       | -                  | -                     | -                | -                        | 1.00E+03           | n            | -                                | 9.70E-03                  | -                        | -                                | 1.09E-01                          | -                                | N                    |
| sec-Butylbenzene               | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Selenium                       | 5.68E+03           | n                     | 1.55E+03         | n                        | 5.10E+03           | n            | 9.65E-01                         | 9.50E-01                  | 2.60E-01                 | 1.09E+01                         | 1.07E+01                          | 2.93E+00                         | N                    |
| Silver                         | 5.68E+03           | n                     | 1.55E+03         | n                        | 5.10E+03           | n            | 1.57E+00                         | 1.60E+00                  | -                        | 1.76E+01                         | 1.80E+01                          | -                                | Y                    |
| Styrene                        | 5.12E+04           | ns                    | 3.03E+04         | ns                       | 3.80E+04           | ns           | 1.56E+00                         | 2.00E+00                  | 1.20E-01                 | 1.76E+01                         | 2.25E+01                          | 1.35E+00                         | N                    |
| 1,2,3-Trichlorobenzene         | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| 1,1,1,2-Tetrachloroethane      | 1.61E+02           | c                     | 2.78E+03         | cs                       | 9.80E+00           | c            | 1.73E-03                         | 2.10E-04                  | -                        | 1.94E-02                         | 2.36E-03                          | -                                | N                    |
| 1,1,1-Trichloroethane          | 7.71E+04           | ns                    | 6.43E+04         | ns                       | 3.90E+04           | ns           | 2.98E+00                         | 3.30E+00                  | 7.20E-02                 | 3.35E+01                         | 3.71E+01                          | 8.10E-01                         | N                    |
| 1,1,2,2-Tetrachloroethane      | 4.33E+01           | c                     | 5.99E+02         | c                        | 2.90E+00           | c            | 2.25E-04                         | 2.80E-05                  | -                        | 2.53E-03                         | 3.15E-04                          | -                                | N                    |
| 1,1,2-Trichloroethane          | 9.43E+01           | c                     | 1.24E+03         | ns                       | 5.50E+00           | c            | 6.74E-04                         | 8.20E-05                  | 1.70E-03                 | 7.58E-03                         | 9.23E-04                          | 1.91E-02                         | N                    |
| 2,4,5-Trichlorophenol          | 6.84E+04           | n                     | 2.38E+04         | n                        | 6.20E+04           | n            | 7.13E+00                         | 9.40E+00                  | -                        | 8.02E+01                         | 1.06E+02                          | -                                | N                    |
| 2,4,6-Trichlorophenol          | 6.84E+02           | n                     | 2.38E+02         | n                        | 1.60E+02           | c**          | 7.13E-02                         | 1.60E-02                  | -                        | 8.02E-01                         | 1.80E-01                          | -                                | N                    |
| 1,2,3-Trichloropropane         | 4.54E+00           | c                     | 3.10E+01         | c                        | 4.10E-01           | c            | 3.56E-05                         | 4.40E-06                  | -                        | 4.01E-04                         | 4.95E-05                          | -                                | N                    |
| 1,2,4-Trichlorobenzene         | 5.25E+02           | ns                    | 4.27E+02         | ns                       | 4.00E+02           | ns           | 1.02E-02                         | 1.30E-02                  | 1.10E-01                 | 1.15E-01                         | 1.46E-01                          | 1.24E+00                         | N                    |
| 1,2,4-Trimethylbenzene         | -                  | -                     | -                | -                        | 2.80E+02           | ns           | -                                | 2.40E-02                  | -                        | -                                | 2.70E-01                          | -                                | N                    |
| 1,3,5-Trimethylbenzene         | -                  | -                     | -                | -                        | 2.00E+02           | n            | -                                | 2.00E-02                  | -                        | -                                | 2.25E-01                          | -                                | N                    |
| tert-Butylbenzene              | -                  | -                     | -                | -                        | -                  | -            | -                                | -                         | -                        | -                                | -                                 | -                                | N                    |
| Tetrachloroethene (PCE)        | 3.64E+01           | c                     | 3.38E+02         | cs                       | 2.70E+00           | c            | 4.49E-04                         | 5.20E-05                  | 2.40E-03                 | 5.05E-03                         | 5.85E-04                          | 2.70E-02                         | N                    |
| Toluene                        | 5.79E+04           | ns                    | 2.11E+04         | ns                       | 4.60E+04           | ns           | 1.38E+00                         | 1.70E+00                  | 7.60E-01                 | 1.56E+01                         | 1.91E+01                          | 8.55E+00                         | Y                    |
| trans-1,2-DCE                  | 9.95E+02           | n                     | 8.14E+02         | n                        | 5.00E+02           | n            | 3.01E-02                         | 3.40E-02                  | 3.20E-02                 | 3.39E-01                         | 3.83E-01                          | 3.60E-01                         | N                    |
| trans-1,3-Dichloropropene      | 1.26E+02           | c                     | 5.10E+02         | n                        | 8.40E+00           | c*           | 1.35E-03                         | 1.60E-04                  | -                        | 1.52E-02                         | 1.80E-03                          | -                                | N                    |
| Trichloroethene (TCE)          | 2.53E+02           | c                     | 4.60E+03         | cs                       | 1.40E+01           | c            | 5.30E-03                         | 6.10E-04                  | 1.90E-03                 | 5.96E-02                         | 6.86E-03                          | 2.14E-02                         | N                    |
| Trichlorofluoromethane         | 6.76E+03           | ns                    | 5.82E+03         | ns                       | 3.40E+03           | ns           | 9.01E-01                         | 8.40E-01                  | -                        | 1.01E+01                         | 9.45E+00                          | -                                | N                    |
| Vanadium                       | 5.68E+03           | n                     | 1.55E+03         | n                        | 7.20E+03           | n            | 1.83E+02                         | 2.60E+02                  | -                        | 2.05E+03                         | 2.93E+03                          | -                                | Y                    |
| Vinyl chloride                 | 2.59E+01           | c                     | 2.48E+02         | c                        | 1.70E+00           | c            | 2.88E-04                         | 5.60E-06                  | 7.00E-04                 | 3.24E-03                         | 6.30E-05                          | 7.88E-03                         | N                    |
| Xylenes, Total                 | 3.61E+03           | ns                    | 3.13E+03         | ns                       | 2.60E+03           | ns           | 1.76E-01                         | 2.30E-01                  | 1.10E+01                 | 1.98E+00                         | 2.59E+00                          | 1.24E+02                         | N                    |
| Zinc                           | 3.41E+05           | nl                    | 9.29E+04         | n                        | 3.10E+05           | nm           | 6.82E+02                         | 6.80E+02                  | -                        | 7.67E+03                         | 7.65E+03                          | -                                | Y                    |

c - carcinogen  
n - noncarcinogen  
cs - carcinogen, SSL may exceed saturation  
ns - noncarcinogen, SSL may exceed saturation  
- no screening value currently available  
nl - noncarcinogen, SSL may exceed ceiling limit  
nls - noncarcinogen, SSL may exceed both saturation and ceiling limit

NMED - Technical Background Document for Development of Soil Screening Levels - Revision 5.0 ( August 2009)  
EPA - Regional Screening Levels (April 2009)

| Analytes   | Sample Date | Residential Soil Screening Level | Leachate DAF (11.25) (mg/kg) SolidW NMED |          | Source | AOC 19-1 (0-0.5') | AOC 19-2 (0-0.5') | AOC 19-3 (0-0.5') | AOC 19-4 (0-0.5') | AOC 19-5 (0-0.5') | AOC 19-6 (0-0.5') | AOC 19-7 (0-0.5') | AOC 19-7 (1.5-2.0) | AOC 19-8 (0-0.5') | AOC 19-9 (0-0.5) | AOC 19-10 (0-0.5') | AOC 19-11 (0-0.5') | AOC 19-12 (0-0.5') | AOC 19-13 (0-0.5') | AOC 19-14 (0-0.5') | AOC 19-15 (0-0.5') | AOC 19-16 (0-0.5') | AOC 19-17 (0-0.5') | AOC 19-18 (0-0.5) | AOC 19-19 (0-0.5') | AOC 19-20 (0-0.5') | AOC 19-21 (0-0.5') | AOC 19-22 (0-0.5') | AOC 19-23 (0-0.5') |     |
|--|-------------|----------------------------------|--|----------|--------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|--------------------|-------------------|------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|-------------------|--------------------|--------------------|--------------------|--------------------|--------------------|-----|
|  |             |                                  | Source                                   | Source   |        |                   |                   |                   |                   |                   |                   |                   |                    |                   |                  |                    |                    |                    |                    |                    |                    |                    |                    |                   |                    |                    |                    |                    |                    |     |
|  |             |                                  | Source                                   | Source   |        |                   |                   |                   |                   |                   |                   |                   |                    |                   |                  |                    |                    |                    |                    |                    |                    |                    |                    |                   |                    |                    |                    |                    |                    |     |
| Metals (mg/kg)                                       |             |                                  |  |          |        |                   |                   |                   |                   |                   |                   |                   |                    |                   |                  |                    |                    |                    |                    |                    |                    |                    |                    |                   |                    |                    |                    |                    |                    |     |
| Antimony   |             | 3.13E+01                         | (1)                                      | 7.44E+00 | (4)    | <2.5              | <12               | <2.5              | <13               | <5.0              | <13               | <2.5              | <2.5               | <2.5              | <2.5             | <5.0               | <12                | <13                | <2.5               | <5.0               | <12                | <2.5               | <2.5               | <2.5              | <2.5               | <2.5               | <2.5               | <2.5               | <2.5               |     |
| Arsenic  |             | 3.59E+00                         | (1)                                      | 1.48E+01 | (4)    | <2.5              | 14                | <2.5              | <13               | <5.0              | <13               | <2.5              | <2.5               | 5.2               | 5.2              | <5.0               | <12                | <13                | 3.6                | <5.0               | <12                | 2.9                | 3.8                | 3.8               | <2.5               | <2.5               | <2.5               | <2.5               | 24                 |     |
| Barium   |             | 1.56E+04                         | (1)                                      | 3.39E+03 | (4)    | 100               | 120               | 23                | 71                | 62                | 27                | 670               | 270                | 96                | 220              | 81                 | 130                | 94                 | 160                | 120                | 230                | 45                 | 120                | 260               | 140                | 100                | 87                 | 110                | 110                |     |
| Beryllium  |             | 1.56E+02                         | (1)                                      | 6.49E+02 | (4)    | 0.26              | 0.88              | 0.28              | 0.84              | 0.31              | 1.4               | 0.51              | 0.43               | 0.35              | 0.26             | <0.30              | <0.75              | 0.87               | 0.45               | 0.51               | 0.93               | 0.21               | 0.19               | 0.45              | 0.34               | 0.17               | 0.15               | 0.21               | <0.15              |     |
| Cadmium  |             | 7.79E+01                         | (1)                                      | 1.55E+01 | (4)    | <0.10             | <0.50             | <0.10             | <0.50             | <0.20             | <0.50             | <0.10             | <0.50              | <0.10             | <0.20            | <0.20              | <0.50              | <0.50              | <0.10              | <0.20              | <0.50              | <0.10              | <0.10              | <0.10             | <0.10              | <0.10              | <0.10              | <0.10              | <0.10              |     |
| Chromium   |             | 2.19E+02                         | (1)                                      | 2.37E+01 | (4)    | 2.8               | 2.4               | 1.6               | 5                 | 2.8               | 8.6               | 3.6               | 2.9                | 3.4               | 2.6              | 1.9                | 20                 | 7                  | 92                 | 600                | 21                 | 34                 | 33                 | 170               | 3.8                | 3.3                | 2.4                | 3.9                | 1.4                |     |
| Cobalt   |             | 2.30E+01                         | (2)                                      | 5.51E+00 | (5)    | 1.6               | 11                | 3.6               | 3.9               | 3.6               | 2.8               | 2.1               | 3.8                | 2.3               | 2.9              | 5.3                | 6                  | 3.1                | 4                  | 7.3                | 2.6                | 1.4                | 2.6                | 2.5               | 2.7                | 1.8                | 2.3                | 2.6                | <0.30              |     |
| Cyanide  |             | 1.56E+03                         | (1)                                      | 8.37E+01 | (4)    | <0.30             | <0.30             | <0.30             | <0.30             | <0.30             | <0.30             | <0.30             | <0.30              | <0.30             | <0.30            | <0.30              | <0.30              | <0.30              | <0.30              | <0.30              | <0.30              | <0.30              | <0.30              | <0.30             | <0.30              | <0.30              | <0.30              | <0.30              | <0.30              |     |
| Lead   |             | 4.00E+02                         | (1)                                      | -        | (4)    | 5.7               | 4.9               | 5.3               | 67                | 30                | 14                | 20                | 6.8                | 6.7               | 3.3              | 2.9                | 11                 | 13                 | 12                 | 10                 | 9.4                | 2.8                | 3.1                | 5.8               | 5.1                | 2.4                | 2.1                | 2.8                | 17                 |     |
| Mercury  |             | 7.71E+00                         | (1)                                      | 3.30E+01 | (4)    | <0.033            | <0.033            | <0.033            | <0.033            | <0.033            | <0.033            | <0.033            | <0.033             | <0.033            | <0.033           | <0.033             | <0.033             | <0.033             | <0.033             | 0.13               | 0.051              | <0.033             | <0.033             | <0.033            | 0.11               | <0.033             | <0.033             | <0.033             | <0.033             |     |
| Nickel   |             | 1.56E+03                         | (1)                                      | 5.36E+02 | (4)    | 2                 | 19                | 3.4               | 5.4               | 3.7               | 6                 | 3.5               | 2.5                | 4                 | 2.5              | 2.8                | 6.3                | 7.4                | 3.8                | 5                  | 9.8                | 2.9                | 1.6                | 3.2               | 3.1                | 3.2                | 3.1                | 2.6                | 1.7                | 2   |
| Selenium   |             | 3.91E+02                         | (1)                                      | 1.09E+01 | (4)    | <2.5              | <12               | <2.5              | <13               | <5.0              | <13               | <2.5              | <2.5               | <2.5              | <2.5             | <5.0               | <12                | <13                | <2.5               | <5.0               | <12                | <2.5               | <2.5               | <2.5              | <2.5               | <2.5               | <2.5               | <2.5               | <2.5               |     |
| Silver   |             | 3.91E+02                         | (1)                                      | 1.76E+01 | (4)    | 0.31              | <1.2              | <0.25             | <1.3              | <0.50             | <1.3              | <0.25             | <0.25              | <0.25             | <0.25            | <0.50              | <1.2               | <1.3               | <0.25              | <0.50              | <1.2               | <0.25              | <0.25              | <0.25             | <0.25              | <0.25              | <0.25              | <0.25              | <0.25              |     |
| Vanadium   |             | 3.91E+02                         | (1)                                      | 2.05E+03 | (4)    | 8.6               | <12               | 6.9               | <12               | 11                | 17                | 7.9               | 9.7                | 11                | 6.6              | <5.0               | 14                 | 15                 | 10                 | 15                 | 28                 | 6.3                | 7.2                | 13                | 12                 | 20                 | 13                 | 13                 | 6.2                | 6.2 |
| Zinc   |             | 2.35E+04                         | (1)                                      | 7.67E+03 | (4)    | 13                | 54                | 17                | 31                | 17                | 43                | 18                | 15                 | 19                | 12               | 16                 | 54                 | 43                 | 110                | 340                | 64                 | 32                 | 43                 | 190               | 20                 | 16                 | 12                 | 15                 | 8.6                |     |
| Volatile Organic Compounds - (EPA Method 8260) mg/kg |             |                                  |  |          |        |                   |                   |                   |                   |                   |                   |                   |                    |                   |                  |                    |                    |                    |                    |                    |                    |                    |                    |                   |                    |                    |                    |                    |                    |     |
| 1,1,1,2-tetrachloroethane                            |             | 2.92E+01                         | (1)                                      | 1.94E+02 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050           | <0.001             | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |     |
| 1,1,1-Trichloroethane                                |             | 2.18E+04                         | (1)                                      | 3.35E+01 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050           | <0.001             | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |     |
| 1,1,2,2-tetrachloroethane                            |             | 7.97E+00                         | (1)                                      | 2.53E+03 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050           | <0.001             | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |     |
| 1,1,2,4-tetrachloroethane                            |             | 6.70E+01                         | (2)                                      | 2.70E+01 | (5)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050           | <0.001             | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |     |
| 1,1,2-Trichloroethane                                |             | 1.72E+01                         | (1)                                      | 7.58E+03 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050           | <0.001             | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |     |
| 1,1,2-Dibromo-3-chloropropene                        |             | 1.94E+01                         | (1)                                      | 3.35E+05 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.10            | <0.001             | <0.001             | <0.001             | <0.10              | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |     |
| 1,2-Dibromoethane (EDB)                              |             | 5.74E+01                         | (1)                                      | 1.78E+04 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050           | <0.001             | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |     |
| 1,2-Dichlorobenzene                                  |             | 3.01E+03                         | (1)                                      | 3.53E+00 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050           | <0.001             | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |     |
| 1,1-Dichloropropene                                  |             | -                                | -  | -        | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.10            | <0.001             | <0.001             | <0.001             | <0.10              | <0.001             | <0.001             | <0.001             | <0.001             |                   |                    |                    |                    |                    |                    |     |

| Analytes   | Sample Date | Residential Soil Screening Level | Source | Leachate DAF (11.25) (mg/kg) Soil/GW NMED | Source |                   |                   |                   |                   |                   |                   |                   |                    |                   |                   |                    |                    |                    |                    |                    |                    |                    |                    |                   |                    |                    |                    |                    |                    |
|--|-------------|----------------------------------|--------|---|--------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|--------------------|-------------------|-------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|-------------------|--------------------|--------------------|--------------------|--------------------|--------------------|
|  |             |                                  |        |   |        | AOC 19-1 (0-0.5') | AOC 19-2 (0-0.5') | AOC 19-3 (0-0.5') | AOC 19-4 (0-0.5') | AOC 19-5 (0-0.5') | AOC 19-6 (0-0.5') | AOC 19-7 (0-0.5') | AOC 19-7 (1.5-2.0) | AOC 19-8 (0-0.5') | AOC 19-9 (0-0.5') | AOC 19-10 (0-0.5') | AOC 19-11 (0-0.5') | AOC 19-12 (0-0.5') | AOC 19-13 (0-0.5') | AOC 19-14 (0-0.5') | AOC 19-15 (0-0.5') | AOC 19-16 (0-0.5') | AOC 19-17 (0-0.5') | AOC 19-18 (0-0.5) | AOC 19-19 (0-0.5') | AOC 19-20 (0-0.5') | AOC 19-21 (0-0.5') | AOC 19-22 (0-0.5') | AOC 19-23 (0-0.5') |
| Methyl tert-butyl ether (MTBE)                   |             | 8.62E+02                         | (1)    | 2.59E-01                                  | (4)    | 0.014             | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |
| Methylene chloride                               |             | 1.99E+02                         | (1)    | 1.21E-01                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | 0.0118            | 0.0125             | <0.001            | <0.15             | <0.001             | <0.001             | <0.001             | <0.15              | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             |
| Naphthalene                                      |             | 4.50E+01                         | (1)    | 4.72E-02                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.10             | <0.001             | <0.001             | <0.10              | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| n-Butylbenzene                                   |             | -                                | -      | -   | -      | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| n-Propylbenzene                                  |             | -                                | -      | -   | -      | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| sec-Butylbenzene                                 |             | -                                | -      | -   | -      | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| Styrene  |             | 8.97E+03                         | (1)    | 1.79E+01                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| tert-Butylbenzene                                |             | -                                | -      | -   | -      | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| Tetrachloroethene (PCE)                          |             | 6.99E+00                         | (1)    | 5.05E-03                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| Toluene  |             | 5.57E+03                         | (1)    | 1.95E+01                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| trans-1,2-DCE                                    |             | 2.73E+02                         | (1)    | 3.39E-01                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| trans-1,3-Dichloropropene                        |             | 2.35E+01                         | (1)    | 1.52E-02                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| Trichloroethene (TCE)                            |             | 4.37E+01                         | (1)    | 5.99E-02                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| Trichloroethenemethane                           |             | 2.01E+03                         | (1)    | 1.01E+01                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| Vinyl chloride                                   |             | 8.65E-01                         | (1)    | 3.24E-03                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.050            | <0.001             | <0.001             | <0.050             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| Xylenes, Total                                   |             | 1.09E+03                         | (1)    | 1.98E+00                                  | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001            | <0.10             | <0.001             | <0.001             | <0.10              | <0.001             | <0.001             | <0.001             | <0.001             | <0.001             | <0.001            | <0.001             | <0.001             | <0.001             | <0.001             |                    |
| Semi Volatile Organics - (EPA Method 8270) mp/kg |             |                                  |        |   |        |                   |                   |                   |                   |                   |                   |                   |                    |                   |                   |                    |                    |                    |                    |                    |                    |                    |                    |                   |                    |                    |                    |                    |                    |
| 1,2,4-Trichlorobenzene                           |             | 1.43E+02                         | (1)    | 1.19E-01                                  | (4)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 1,2-Dichlorobenzene                              |             | 3.01E+03                         | (1)    | 3.53E+00                                  | (4)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 1,3-Dichlorobenzene                              |             | -                                | -      | -   | -      | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 1,4-Dichlorobenzene                              |             | 3.21E+01                         | (1)    | 4.02E-02                                  | (4)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 1-Methylnaphthalene                              |             | 2.20E+02                         | (1)    | 1.69E-01                                  | (5)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 2,4,5-Trichlorophenol                            |             | 6.11E+03                         | (1)    | 8.02E+01                                  | (4)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 2,4,6-Trichlorophenol                            |             | 6.11E+01                         | (1)    | 8.02E-01                                  | (4)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 2,4-Dichlorophenol                               |             | 1.83E+02                         | (1)    | 1.54E+00                                  | (4)    | <0.40             | <0.40             | <0.40             | <0.40             | <0.40             | <0.40             | <0.80             | <0.40              | <0.40             | <0.40             | <0.40              | <2.0               | <0.40              | <0.80              | <2.0               | <0.80              | <0.80              | <0.80              | <0.40             | <0.80              | <0.40              | <0.40              | <0.80              | <2.0               |
| 2,4-Dimethylphenol                               |             | 1.22E+03                         | (1)    | 1.03E+01                                  | (4)    | <0.30             | <0.30             | <0.30             | <0.30             | <0.30             | <0.30             | <0.60             | <0.30              | <0.30             | <0.30             | <0.30              | <1.5               | <0.30              | <0.60              | <1.5               | <0.60              | <0.60              | <0.60              | <0.30             | <0.60              | <0.30              | <0.30              | <0.60              | <1.5               |
| 2,4-Dinitrophenol                                |             | 1.22E+02                         | (1)    | 5.91E-01                                  | (4)    | <0.40             | <0.40             | <0.40             | <0.40             | <0.40             | <0.40             | <0.80             | <0.40              | <0.40             | <0.40             | <0.40              | <2.0               | <0.40              | <0.80              | <2.0               | <0.80              | <0.80              | <0.80              | <0.40             | <0.80              | <0.40              | <0.40              | <0.80              | <2.0               |
| 2,6-Dinitrophenol                                |             | 1.26E+01                         | (1)    | 1.75E-02                                  | (4)    | <0.50             | <0.50             | <0.50             | <0.50             | <0.50             | <0.50             | <1.0              | <0.50              | <0.50             | <0.50             | <0.50              | <2.5               | <0.50              | <1.0               | <2.5               | <1.0               | <1.0               | <1.0               | <0.50             | <1.0               | <0.50              | <0.50              | <1.0               | <2.5               |
| 2,6-Dinitrotoluene                               |             | 6.12E+01                         | (1)    | 3.00E-01                                  | (4)    | <0.50             | <0.50             | <0.50             | <0.50             | <0.50             | <0.50             | <1.0              | <0.50              | <0.50             | <0.50             | <0.50              | <2.5               | <0.50              | <1.0               | <2.5               | <1.0               | <1.0               | <1.0               | <0.50             | <1.0               | <0.50              | <0.50              | <1.0               | <2.5               |
| 2-Chloronaphthalene                              |             | 6.26E+03                         | (1)    | 1.52E+02                                  | (4)    | <0.25             | <0.25             | <0.25             | <0.25             | <0.25             | <0.25             | <0.50             | <0.25              | <0.25             | <0.25             | <0.25              | <1.2               | <0.25              | <0.50              | <1.2               | <0.50              | <0.50              | <0.50              | <0.25             | <0.50              | <0.25              | <0.25              | <0.50              | <1.2               |
| 2-Chlorophenol                                   |             | 3.91E+02                         | (1)    | 1.72E+00                                  | (4)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 2-Methylnaphthalene                              |             | 3.10E+02                         | (2)    | 1.01E+01                                  | (5)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 2-Methylphenol                                   |             | 3.10E+03                         | (2)    | 2.25E+01                                  | (5)    | <0.50             | <0.50             | <0.50             | <0.50             | <0.50             | <0.50             | <1.0              | <0.50              | <0.50             | <0.50             | <0.50              | <2.5               | <0.50              | <1.0               | <2.5               | <1.0               | <1.0               | <1.0               | <0.50             | <1.0               | <0.50              | <0.50              | <1.0               | <2.5               |
| 2-Nitroaniline                                   |             | 1.80E+02                         | (2)    | 3.71E-01                                  | (5)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 2-Nitrophenol                                    |             | -                                | -      | -   | -      | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | <0.20              | <0.20             | <0.20             | <0.20              | <1.0               | <0.20              | <0.40              | <1.0               | <0.40              | <0.40              | <0.40              | <0.20             | <0.40              | <0.20              | <0.20              | <0.40              | <1.0               |
| 3,3-Dichlorobenzidine                            |             | 8.71E+00                         | (1)    | 1.92E-01                                  | (4)    | <0.25             | <0.25             | <0.25             | <0.25             | <0.25             | <0.25             | <0.50             | <0.25              | <0.25             | <0.25             | <0.25              | <1.2               | <0.25              | <0.50              | <1.2               | <0.50              | <0.50              | <0.50              | <0.25             | <0.50              | <0.25              | <0.25              | <0.50              | <1.2               |
| 3,4-Methylphenol                                 |             | 3.10E+02                         | (2)    | 2.14E+00                                  | (5)    | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20             | <0.40             | &                  |                   |                   |                    |                    |                    |                    |                    |                    |                    |                    |                   |                    |                    |                    |                    |                    |



| Analyses       | Sample Date | Residential Soil Screening Level | Leachate DAF (11.25) Soil/GW NMED |        | AOC               |           |                   |           |                   |           |                   |           |                   |           |                   |           |                   |           |                   |           |                   |           |                    |           |                    |           |                    |           |                     |           |                     |           |                   |           |                   |           |                   |           |                   |           |                   |           |                     |           |                     |           |           |           |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
|----------------|-------------|----------------------------------|-----------------------------------|--------|-------------------|-----------|-------------------|-----------|-------------------|-----------|-------------------|-----------|-------------------|-----------|-------------------|-----------|-------------------|-----------|-------------------|-----------|-------------------|-----------|--------------------|-----------|--------------------|-----------|--------------------|-----------|---------------------|-----------|---------------------|-----------|-------------------|-----------|-------------------|-----------|-------------------|-----------|-------------------|-----------|-------------------|-----------|---------------------|-----------|---------------------|-----------|-----------|-----------|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|--|
|                |             |                                  | Source                            | Source | AOC 20-1 (0-0.5') |           | AOC 20-2 (0-0.5') |           | AOC 20-3 (0-0.5') |           | AOC 20-4 (0-0.5') |           | AOC 20-5 (0-0.5') |           | AOC 20-6 (0-0.5') |           | AOC 20-7 (0-0.5') |           | AOC 20-8 (0-0.5') |           | AOC 20-9 (0-0.5') |           | AOC 20-10 (0-0.5') |           | AOC 20-11 (0-0.5') |           | AOC 20-12 (0-0.5') |           | AOC 21-1 (1.5-2.0') |           | AOC 21-1 (2.0-2.5') |           | AOC 21-1-(0-0.5') |           | AOC 21-2 (0-0.5') |           | AOC 21-3 (0-0.5') |           | AOC 21-4 (0-0.5') |           | AOC 21-5 (0-0.5') |           | AOC 21-5 (1.5-2.0') |           | AOC 21-5 (2.0-2.5') |           |           |           |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
|                |             |                                  |                                   |        | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011          | 8/17/2011 | 8/17/2011          | 8/17/2011 | 8/17/2011          | 8/17/2011 | 8/17/2011           | 8/17/2011 | 8/17/2011           | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011         | 8/17/2011 | 8/17/2011           | 8/17/2011 | 8/17/2011           | 8/17/2011 | 8/17/2011 | 8/17/2011 |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| Metals (mg/kg) |             |                                  |                                   |        |                   |           |                   |           |                   |           |                   |           |                   |           |                   |           |                   |           |                   |           |                   |           |                    |           |                    |           |                    |           |                     |           |                     |           |                   |           |                   |           |                   |           |                   |           |                   |           |                     |           |                     |           |           |           |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |  |



| Analytes   | Sample Date | Residential Soil Screening Level | Leachate DAF (11.25) (mg/kg) Soil/GW NMED |          | Source | AOC 20-1 (0-0.5') | AOC 20-2 (0-0.5') | AOC 20-3 (0-0.5') | AOC 20-4 (0-0.5') | AOC 20-5 (0-0.5') | AOC 20-6 (0-0.5') | AOC 20-7 (0-0.5') | AOC 20-8 (0-0.5') | AOC 20-9 (0-0.5') | AOC 20-10 (0-0.5') | AOC 20-11 (0-0.5') | AOC 20-12 (0-0.5') | AOC 21-1 (1.5-2.0') | AOC 21-1 (2.0-2.5') | AOC 21-1(0-0.5') | AOC 21-2 (0-0.5') | AOC 21-3 (0-0.5') | AOC 21-4 (0-0.5') | AOC 21-5 (0-0.5') | AOC 21-5 (1.5-2.0') | AOC 21-5 (2.0-2.5') |
|--|-------------|----------------------------------|---|----------|--------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|--------------------|--------------------|--------------------|---------------------|---------------------|------------------|-------------------|-------------------|-------------------|-------------------|---------------------|---------------------|
|  |             |                                  | Source                                    | Source   |        |                   |                   |                   |                   |                   |                   |                   |                   |                   |                    |                    |                    |                     |                     |                  |                   |                   |                   |                   |                     |                     |
|  |             |                                  | Source                                    | Source   |        |                   |                   |                   |                   |                   |                   |                   |                   |                   |                    |                    |                    |                     |                     |                  |                   |                   |                   |                   |                     |                     |
| Methyl tert-butyl ether (MTBE)                   |             | 8.62E+02                         | (1)                                       | 2.58E-01 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Methylene chloride                               |             | 1.99E+02                         | (1)                                       | 1.21E-01 | (4)    | 0.00809           | 0.0117            | 0.0148            | 0.00884           | 0.00975           | <0.001            | <0.001            | 0.0044            | <0.001            | <0.001             | <0.001             | <0.001             | 0.00275             | 0.00369             | 0.00398          | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Naphthalene                                      |             | 4.50E+01                         | (1)                                       | 4.72E-02 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| n-Butylbenzene                                   |             | -                                | -   | -        | -      | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| n-Propylbenzene                                  |             | -                                | -   | -        | -      | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| sec-Butylbenzene                                 |             | -                                | -   | -        | -      | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Styrene  |             | 8.97E+03                         | (1)                                       | 1.76E+01 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| tert-Butylbenzene                                |             | -                                | -   | -        | -      | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Tetrachloroethene (PCE)                          |             | 6.99E+03                         | (1)                                       | 5.05E-03 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Toluene  |             | 5.57E+03                         | (1)                                       | 1.56E+01 | (4)    | 0.00131           | <0.001            | 0.0143            | 0.00566           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| trans-1,2-DCP                                    |             | 2.73E+02                         | (1)                                       | 3.39E-01 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| trans-1,3-Dichloropropene                        |             | 2.35E+01                         | (1)                                       | 1.52E-02 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Trichloroethene (TCE)                            |             | 4.57E+01                         | (1)                                       | 5.96E-02 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Trichlorofluoromethane                           |             | 2.01E+03                         | (1)                                       | 1.01E+01 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Vinyl chloride                                   |             | 8.65E-01                         | (1)                                       | 3.24E-03 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Xylenes, Total                                   |             | 1.09E+03                         | (1)                                       | 1.98E+00 | (4)    | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001            | <0.001             | <0.001             | <0.001             | <0.001              | <0.001              | <0.001           | <0.001            | <0.001            | <0.001            | <0.001            | <0.001              | <0.001              |
| Semi Volatile Organics - (EPA Method 8270) mg/kg |             |                                  |   |          | (4)    |                   |                   |                   |                   |                   |                   |                   |                   |                   |                    |                    |                    |                     |                     |                  |                   |                   |                   |                   |                     |                     |
| 1,2,4-Trichlorobenzene                           |             | 1.43E+02                         | (1)                                       | 1.15E-01 | (4)    | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 1,2-Dichlorobenzene                              |             | 3.01E+03                         | (1)                                       | 3.53E+00 | (4)    | <0.20             | <0.20             | <0.20             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 1,3-Dichlorobenzene                              |             | -                                | -   | -        | -      | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 1,4-Dichlorobenzene                              |             | 3.21E+01                         | (1)                                       | 4.02E-02 | (4)    | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 1-Methylnaphthalene                              |             | 2.20E+02                         | (3)                                       | 1.69E-01 | (5)    | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 2,4,5-Trichlorophenol                            |             | 6.11E+01                         | (1)                                       | 8.02E+01 | (4)    | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 2,4,6-Trichlorophenol                            |             | 8.20E+01                         | (1)                                       | 8.02E-01 | (4)    | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 2,4-Dichlorophenol                               |             | 1.83E+02                         | (1)                                       | 1.54E+00 | (4)    | <0.40             | <0.40             | <0.80             | <0.80             | <0.80             | <0.30             | <0.30             | <0.30             | <0.40             | <0.40              | <0.40              | <0.40              | <0.40               | <0.40               | <0.40            | <0.40             | <0.40             | <0.40             | <0.40             | <0.40               | <0.40               |
| 2,4-Dimethylphenol                               |             | 1.22E+03                         | (1)                                       | 1.03E+01 | (4)    | <0.30             | <0.30             | <0.60             | <0.60             | <0.60             | <0.30             | <0.30             | <0.30             | <0.30             | <0.30              | <0.30              | <0.30              | <0.30               | <0.30               | <0.30            | <0.30             | <0.30             | <0.30             | <0.30             | <0.30               | <0.30               |
| 2,4-Dinitrophenol                                |             | 1.22E+02                         | (1)                                       | 5.91E-01 | (4)    | <0.40             | <0.40             | <0.80             | <0.80             | <0.80             | <0.40             | <0.40             | <0.40             | <0.40             | <0.40              | <0.40              | <0.40              | <0.40               | <0.40               | <0.40            | <0.40             | <0.40             | <0.40             | <0.40             | <0.40               | <0.40               |
| 2,4-Dinitrotoluene                               |             | 1.26E+01                         | (1)                                       | 1.75E-02 | (4)    | <0.50             | <0.50             | <0.99             | <0.99             | <1.0              | <0.50             | <0.50             | <0.50             | <0.50             | <0.50              | <0.50              | <0.50              | <0.50               | <0.50               | <0.50            | <0.50             | <0.50             | <0.50             | <0.50             | <0.50               | <0.50               |
| 2,6-Dinitrotoluene                               |             | 6.12E+01                         | (1)                                       | 3.00E-01 | (4)    | <0.50             | <0.50             | <0.99             | <1.0              | <1.0              | <0.50             | <0.50             | <0.50             | <0.50             | <0.50              | <0.50              | <0.50              | <0.50               | <0.50               | <0.50            | <0.50             | <0.50             | <0.50             | <0.50             | <0.50               | <0.50               |
| 2-Chloronaphthalene                              |             | 6.26E+03                         | (1)                                       | 1.52E+02 | (4)    | <0.25             | <0.25             | <0.40             | <0.40             | <0.40             | <0.25             | <0.25             | <0.25             | <0.25             | <0.25              | <0.25              | <0.25              | <0.25               | <0.25               | <0.25            | <0.25             | <0.25             | <0.25             | <0.25             | <0.25               | <0.25               |
| 2-Chlorophenol                                   |             | 3.91E+02                         | (1)                                       | 1.72E+00 | (4)    | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 2-Methylnaphthalene                              |             | 3.10E+02                         | (2)                                       | 1.01E+01 | (5)    | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 2-Methylphenol                                   |             | 3.10E+03                         | (2)                                       | 2.25E+01 | (5)    | <0.50             | <0.50             | <0.99             | <1.0              | <1.0              | <0.50             | <0.50             | <0.50             | <0.50             | <0.50              | <0.50              | <0.50              | <0.50               | <0.50               | <0.50            | <0.50             | <0.50             | <0.50             | <0.50             | <0.50               | <0.50               |
| 2-Nitroaniline                                   |             | 1.80E+02                         | (2)                                       | 3.71E-01 | (5)    | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 2-Nitrophenol                                    |             | -                                | -   | -        | -      | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 3,3'-Dichlorobenzidine                           |             | 8.71E+00                         | (1)                                       | 1.92E-01 | (4)    | <0.25             | <0.25             | <0.50             | <0.50             | <0.50             | <0.25             | <0.25             | <0.25             | <0.25             | <0.25              | <0.25              | <0.25              | <0.25               | <0.25               | <0.25            | <0.25             | <0.25             | <0.25             | <0.25             | <0.25               | <0.25               |
| 3+4-Methylphenol                                 |             | 3.10E+02                         | (2)                                       | 2.14E+00 | (5)    | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 3-Nitroaniline                                   |             | -                                | -   | -        | -      | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 4,6-Dinitro-2-methylphenol                       |             | -                                | -   | -        | -      | <0.50             | <0.50             | <0.99             | <1.0              | <1.0              | <0.50             | <0.50             | <0.50             | <0.50             | <0.50              | <0.50              | <0.50              | <0.50               | <0.50               | <0.50            | <0.50             | <0.50             | <0.50             | <0.50             | <0.50               | <0.50               |
| 4-Chloro-3-methylphenol                          |             | -                                | -   | -        | -      | <0.20             | <0.20             | <0.40             | <0.40             | <0.40             | <0.20             | <0.20             | <0.20             | <0.20             | <0.20              | <0.20              | <0.20              | <0.20               | <0.20               | <0.20            | <0.20             | <0.20             | <0.20             | <0.20             | <0.20               | <0.20               |
| 4-Chloroaniline                                  |             | 2.40E+01                         | (3)                                       | 1.35E-03 | (5)    | <0.50             | <0.50             | <0.99             | <1.0              | <1.0              | <0.50             | <0.50             | <0.50             | <0.50             | <0.50              | <0.50              | <0.50              | <0.50               | <0.50               | <0.50            | <0.50             | <0.50             | <0.50             | <                 |                     |                     |

| Analytes          | Sample Date | Residential Soil Screening Level | Source | Leachate DAF (11.25) (mg/kg) Soil/GW NMED | Source | Sample Date |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |     |
|-------------------|-------------|----------------------------------|--------|---|--------|-------------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-----|
|                   |             |                                  |        |   |        |             |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |     |
|                   |             |                                  |        |   |        |             |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |       |     |
| Diethyl phthalate |             | 4.89E+04                         | (1)    | 1.19E+02                                  | (4)    | <0.20       | <0.20 | <0.40 | <0.40 | <0.40 | <0.20 | <0.20 | <0.40 | <0.40 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | <0.20 | &lt |

– No screening level or analytical result available

NMED - Technical Background Document for Development of Soil Screening Levels - Revision 5.0 ( August 2009)

EPA - Regional Screening Levels (April 2009)

(1) NMED Residential Screening Level

(2) EPA Residential Screening Level

(3) EPA Residential - Screening Levels (April 2009) multiplied by 10 pursuant to Provision VII.B. of the July 7, 2007 NMED Order because the constituent is listed as carcinogenic

(4) SoilGW NMED Dilution Attenuation Factor (DAF) = 11.25

(5) SoilGW Risk-based EPA DAF = 11.25

(6) NMED Oct. 2006 TPH Screening Guidelines "unknown oil" with DAF = 1.0 - see report Section 5 for use of TPH screening levels

(7) NMED Oct. 2006 TPH Screening Guidelines(updated with Massachusetts Department of Environmental Protection 2009 tox data - unknown oil, residential, direct contact pathways

(8) NMED Oct. 2006 TPH Screening Guidelines "waste oil" - see report Section 5 for use of TPH screening levels

(9) NMED Oct. 2006 TPH Screening Guidelines "waste oil" with DAF = 1.0 - see report Section 5 for use of TPH screening levels

**Bold represents value above Residential Screening Level**  
 yellow highlight represents value above Leachate (DAF) Screening Level  
**Bold with yellow highlight value exceeds Residential Screening Level and DAF**

**Table 6**  
**Group 6 Soil Cumulative Risk and Hazard Index Evaluation**  
**Bloomfield Refinery - Bloomfield, New Mexico**

| Analytes  | Residential Soil Screening Level | Source | Non-Residential Soil Screening Level | Source | Maximum Concentration | Maximum Concentration divided by Residential Soil Screening Level | Maximum Concentration divided by Non-Residential Soil Screening Level |
|---|----------------------------------|--------|--------------------------------------|--------|-----------------------|---|---|
| <b>CARCINOGENIC CONSTITUENTS</b>                                  |                                  |        |                                      |        |                       |   |   |
| <b>Metals (mg/kg)</b>   |                                  |        |                                      |        |                       |   |   |
| Nickel  | 1.56E+03                         | (1)    | 6.19E+03                             | (5)    | 19                    | 0.0121  | 0.0031  |
| Arsenic   | 3.59E+00                         | (1)    | 1.77E+01                             | (4)    | 24                    | 6.6907  | 1.3565  |
| <b>Volatile Organic Compounds - (EPA Method 8260) mg/kg</b>       |                                  |        |                                      |        |                       |   |   |
| Methylene chloride  | 1.99E+02                         | (1)    | 1.09E+03                             | (4)    | 0.0149                | 0.00007   | 0.00001   |
| <b>Carcinogenic Constituents Cumulative Risk x 10<sup>5</sup></b> |                                  |        |                                      |        |                       | <b>6.703</b>  | <b>1.360</b>  |
| <b>NON-CARCINOGENIC CONSTITUENTS</b>                              |                                  |        |                                      |        |                       |   |   |
| <b>Metals (mg/kg)</b>   |                                  |        |                                      |        |                       |   |   |
| Barium  | 1.56E+04                         | (1)    | 4.35E+03                             | (5)    | 670                   | 0.0430  | 0.1539  |
| Beryllium   | 1.56E+02                         | (1)    | 1.44E+02                             | (5)    | 2.1                   | 0.0134  | 0.0146  |
| Chromium  | 2.19E+02                         | (1)    | 4.49E+02                             | (5)    | 600                   | 2.7397  | 1.3363  |
| Cobalt  | 2.30E+01                         | (2)    | 3.00E+02                             | (6)    | 11                    | 0.4783  | 0.0367  |
| Silver  | 3.91E+02                         | (1)    | 1.55E+03                             | (5)    | 0.31                  | 0.0008  | 0.0002  |
| Vanadium  | 3.91E+02                         | (1)    | 1.55E+03                             | (5)    | 32                    | 0.0818  | 0.0207  |
| Zinc  | 2.35E+04                         | (1)    | 9.29E+04                             | (5)    | 340                   | 0.0145  | 0.0037  |
| Lead  | 4.00E+02                         | (1)    | 8.00E+02                             | (4)    | 67                    | 0.1675  | 0.0838  |
| Mercury   | 7.71E+00                         | (1)    | 4.99E+01                             | (4)    | 0.13                  | 0.0169  | 0.0026  |
| <b>Volatile Organic Compounds - (EPA Method 8260) mg/kg</b>       |                                  |        |                                      |        |                       |   |   |
| 2-Butanone (MEK)  | 3.96E+04                         | (1)    | 1.48E+05                             | (5)    | 0.0438                | 0.000001  | 0.0000003   |
| 2-Hexanone  | -                                | -      | -                                    | -      | 0.00584               | -   | -   |
| 4-Methyl-2-pentanone  | -                                | -      | -                                    | -      | 0.00244               | -   | -   |
| Acetone   | 6.75E+04                         | (1)    | 2.63E+05                             | (5)    | 0.134                 | 0.000002  | 0.000001  |
| 4-Isopropyltoluene  | -                                | -      | -                                    | -      | 0.00477               | -   | -   |
| Carbon disulfide  | 1.94E+03                         | (1)    | 5.89E+03                             | (5)    | 0.00461               | 0.000002  | 0.000001  |
| Toluene   | 5.57E+03                         | (1)    | 2.11E+04                             | (5)    | 0.0256                | 0.000005  | 0.000001  |
| <b>Semi Volatile Organics - (EPA Method 8270) mg/kg</b>           |                                  |        |                                      |        |                       |   |   |
| Benzoic acid  | 2.40E+05                         | (2)    | 2.50E+06                             | (6)    | 3.4                   | 0.0000  | 0.0000  |
| 3+4-Methylphenol  | 3.10E+02                         | (2)    | 3.10E+03                             | (6)    | 0.27                  | 0.0009  | 0.0001  |
| <b>Non-Carcinogenic Constituents Hazard Index</b>                 |                                  |        |                                      |        |                       | <b>3.557</b>  | <b>1.652</b>  |

- No screening level or analytical result available

NMED - Technical Background Document for Development of Soil Screening Levels - Revision 5.0 ( August 2009)

EPA - Regional Screening Levels (April 2009)

(1) NMED Residential Screening Level

(2) EPA Residential Screening Level

(3) EPA Residential - Screening Levels (April 2009) multiplied by 10 pursuant to Provision VII.B. of the July 7, 2007 NMED Order because the constituent is listed as carcinogenic

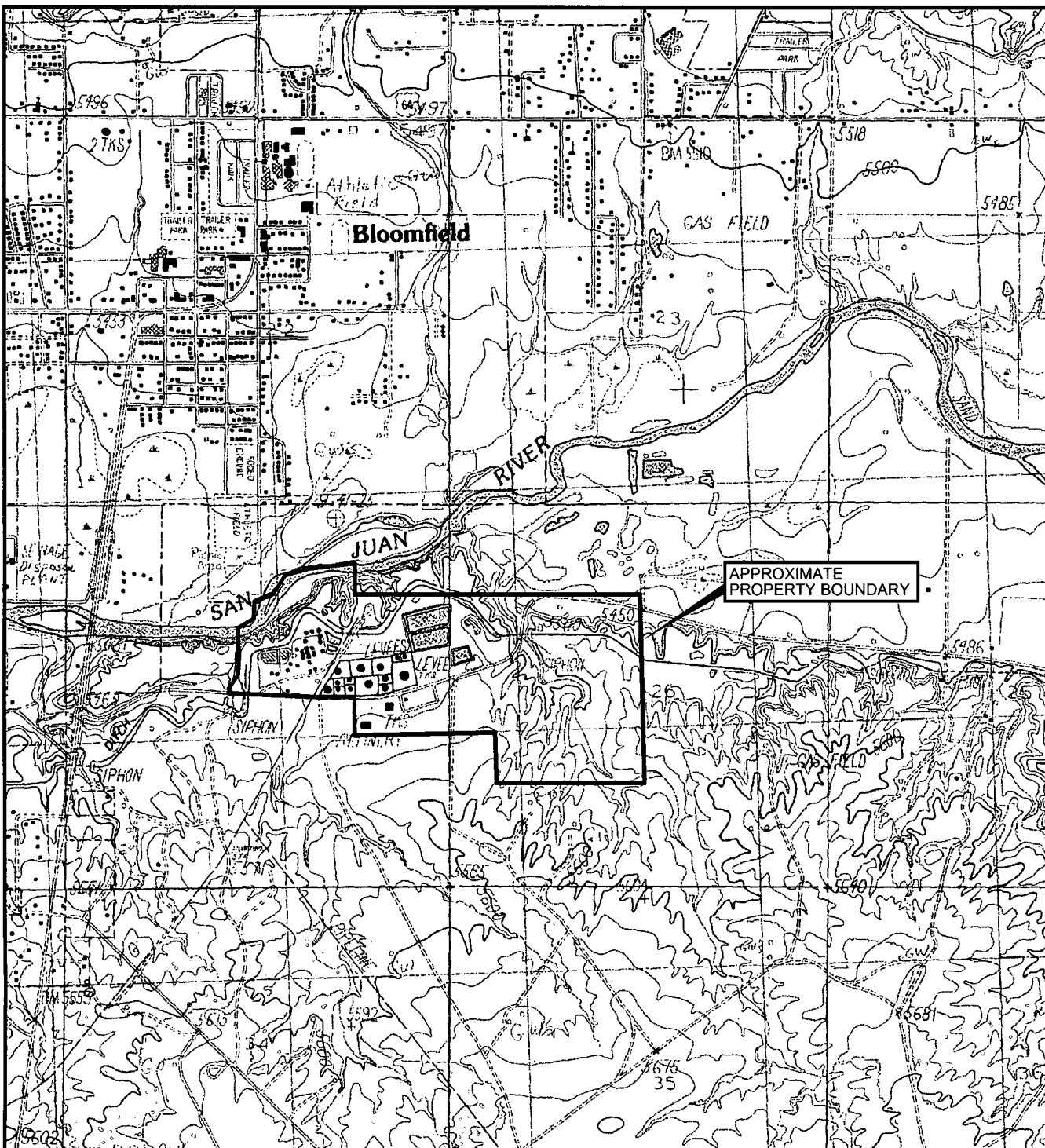
(4) NMED Industrial Occupational Screening Level

(5) NMED Construction Worker Screening Level

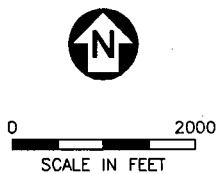
(6) EPA Industrial - Screening Levels (April 2009)

(7) EPA Industrial - Screening Levels (April 2009) multiplied by 10 pursuant to Provision VII.B. of the July 7, 2007 NMED Order because the constituent is listed as carcinogenic





Map Source: USGS 7.5 Min. Quad Sheet BLOOMFIELD, NM., 1985.



**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 01/06/10 FILE: WestRef-A25

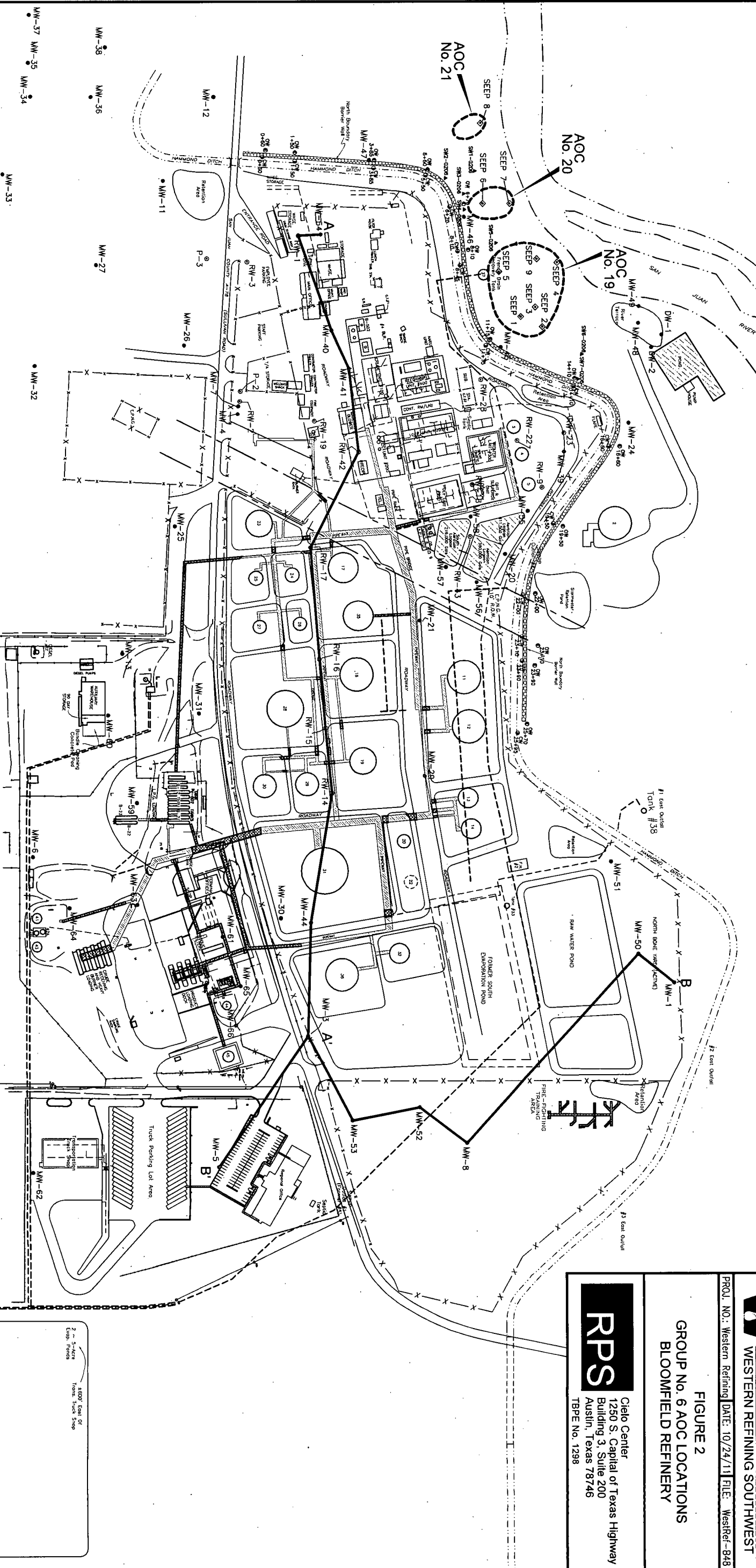
FIGURE 1  
SITE LOCATION MAP  
BLOOMFIELD REFINERY

**RPS**

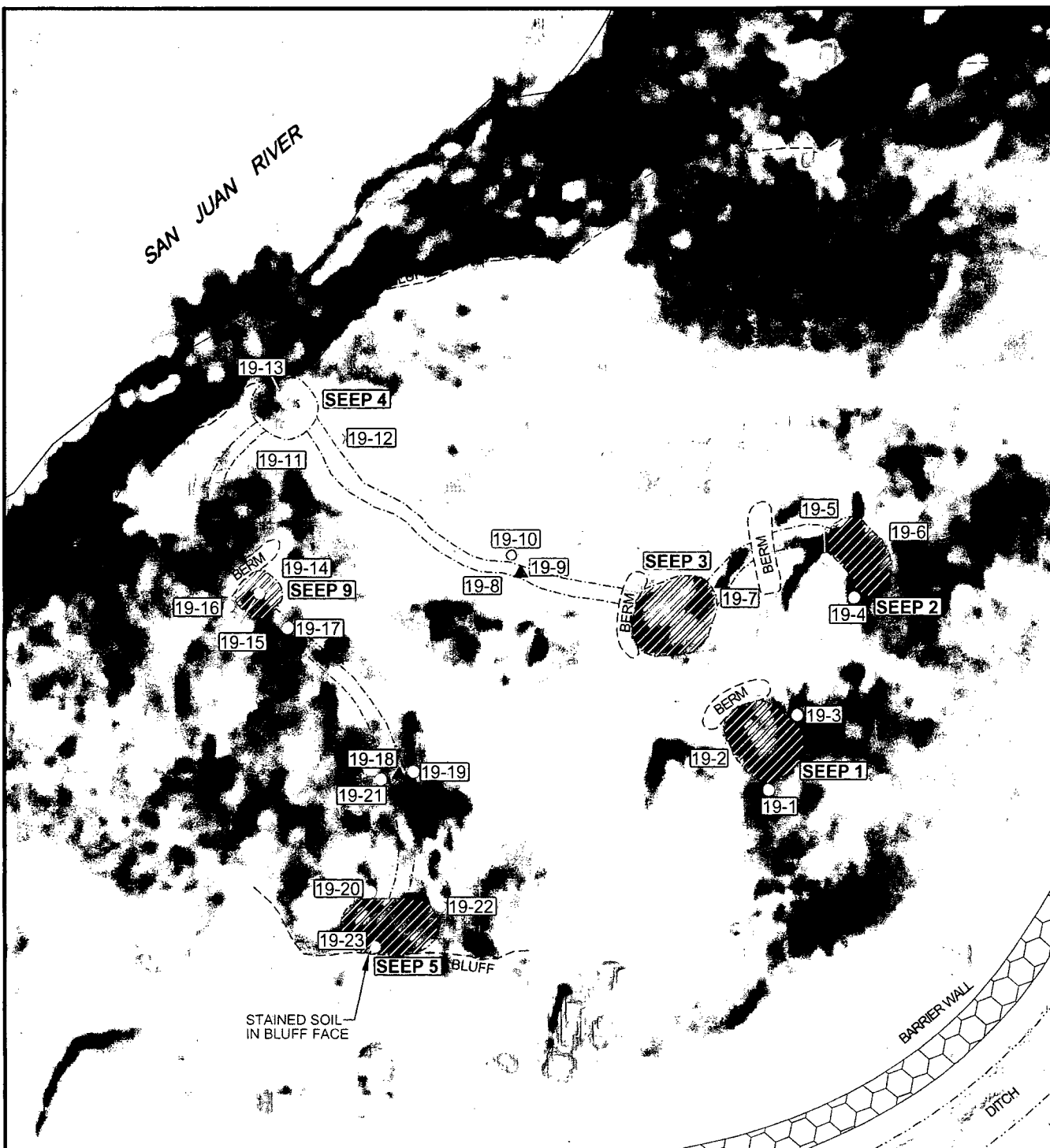
404 Camp Craft Road  
Austin, Texas 78746

**FIGURE 2**  
**GROUP No. 6 AOC LOCATIONS**  
**BLOOMFIELD REFINERY**

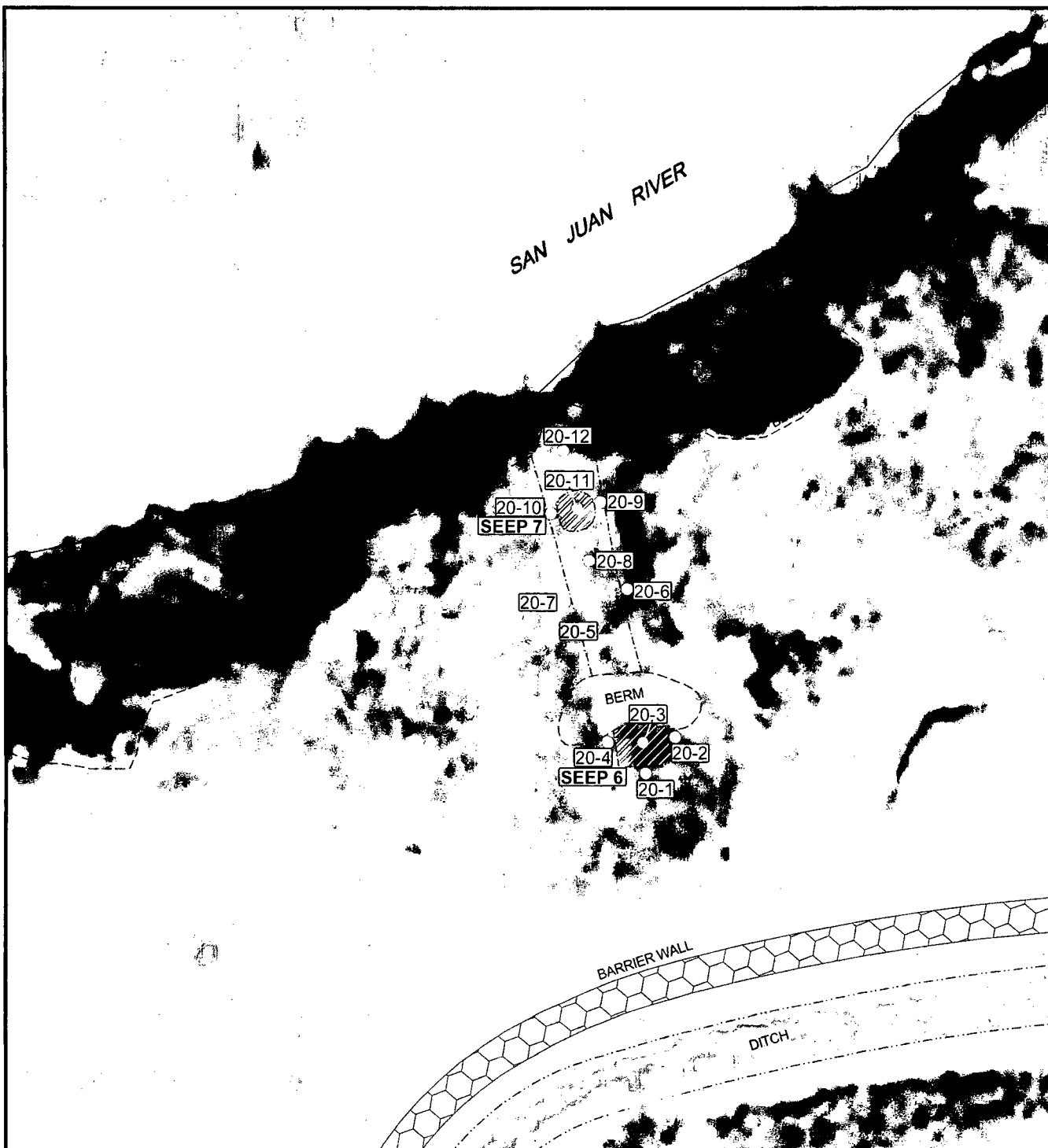
**RPS**  
Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298








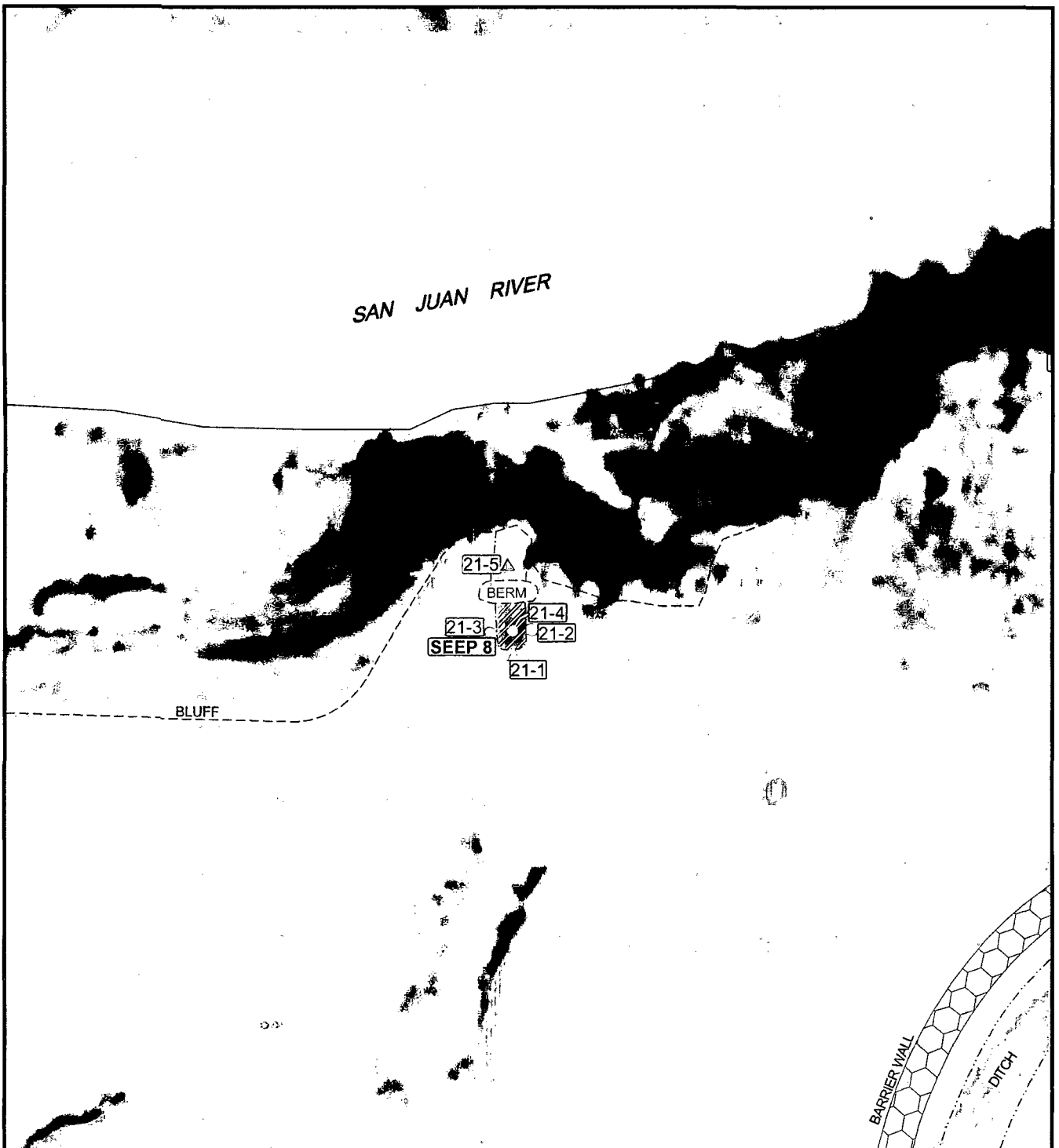




|  |  |  |
|--|--|--|
| <div style="text-align: center;"> </div> <div style="text-align: center; margin-top: 10px;"> <p>SCALE IN FEET</p> </div> | <p style="text-align: center;"><b>LEGEND</b></p> <p><b>SEEP 1</b> SEEP IDENTIFICATION NUMBER</p> <p>○ SURFACE SOIL SAMPLE LOCATION</p> <p>△ SOIL BORING LOCATION</p> <p> LINED CATCHMENT BASIN</p> <p>--- DRAINAGE PATHWAY</p> | <div style="text-align: center;"> <p><b>Western Refining</b><br/>WESTERN REFINING SOUTHWEST</p> </div> <div style="border: 1px solid black; padding: 2px; font-size: 0.8em; margin-top: 5px;">             PROJ. NO.: Western Refining   DATE: 10/24/11   FILE: WestRef-A135         </div> <div style="text-align: center; margin-top: 10px;"> <p><b>FIGURE 4</b><br/><b>AOC No. 19</b><br/><b>SAMPLE LOCATION MAP</b><br/><b>BLOOMFIELD REFINERY</b></p> </div> <div style="display: flex; align-items: center; margin-top: 10px;"> <div style="border: 1px solid black; padding: 5px; text-align: center; flex: 1;"> </div> <div style="flex: 1; padding-left: 10px; font-size: 0.8em;"> <p>Cielo Center<br/>1250 S. Capital of Texas Highway<br/>Building 3, Suite 200<br/>Austin, Texas 78746<br/>TBPE No. 1298</p> </div> </div> |
|--|--|--|



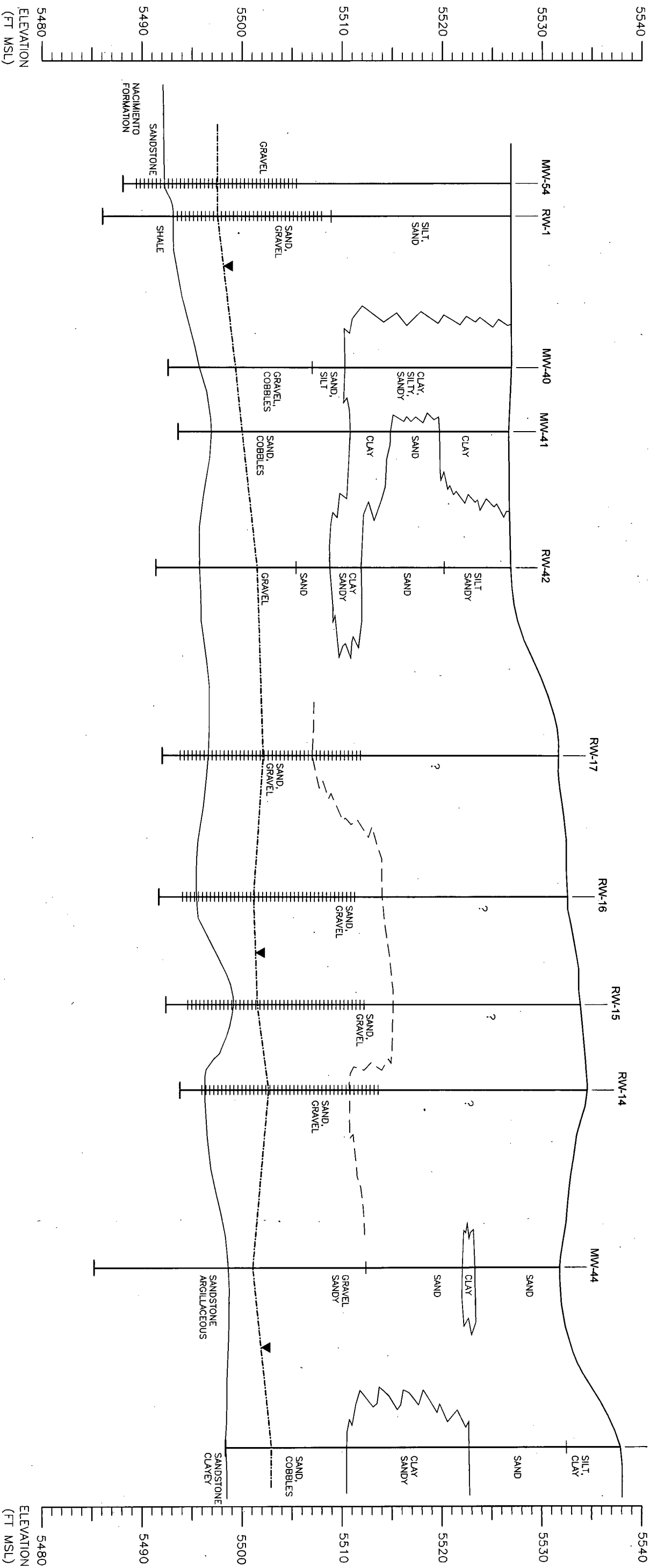
|   |   |
|---|---|
| <div style="text-align: center;">  </div> <div style="margin-top: 10px;">  <p>SCALE IN FEET</p> </div> <div style="margin-top: 20px;"> <p style="text-align: center;"><u>LEGEND</u></p> <p><b>SEEP 6</b> SEEP IDENTIFICATION NUMBER</p> <p>○ SURFACE SOIL SAMPLE LOCATION</p> <p>△ SOIL BORING LOCATION</p> <p> LINED CATCHMENT BASIN</p> <p>----- DRAINAGE PATHWAY</p> </div> | <div style="text-align: center;">  <p><b>Western Refining</b><br/>WESTERN REFINING SOUTHWEST</p> </div> <div style="border: 1px solid black; padding: 5px; margin-top: 5px;"> <p>PROJ. NO.: Western Refining   DATE: 10/24/11   FILE: WestRef-A135</p> </div> <div style="text-align: center; margin-top: 10px;"> <p><b>FIGURE 5</b><br/><b>AOC No. 20</b><br/><b>SAMPLE LOCATION MAP</b><br/><b>BLOOMFIELD REFINERY</b></p> </div> <div style="display: flex; justify-content: space-between; align-items: center; margin-top: 10px;"> <div style="border: 1px solid black; padding: 5px; text-align: center;">  </div> <div style="text-align: right;"> <p>Cielo Center<br/>1250 S. Capital of Texas Highway<br/>Building 3, Suite 200<br/>Austin, Texas 78746<br/>TBPE No. 1298</p> </div> </div> |
|---|---|



|   |  |   |  |
|---|--|---|--|
|   |  | <b>Western Refining</b><br>WESTERN REFINING SOUTHWEST   |  |
| 0  50<br>SCALE IN FEET  |  | PROJ. NO.: Western Refining   DATE: 10/24/11   FILE: WestRef-A135   |  |
| <b>LEGEND</b><br>SEEP 8 SEEP IDENTIFICATION NUMBER<br>○ SURFACE SOIL SAMPLE LOCATION<br>Δ SOIL BORING LOCATION<br>LINED CATCHMENT BASIN<br>- - - DRAINAGE PATHWAY |  | <b>FIGURE 6</b><br>AOC No. 21<br>SAMPLE LOCATION MAP<br>BLOOMFIELD REFINERY                                       |  |
|   |  | Cielo Center<br>1250 S. Capital of Texas Highway<br>Building 3, Suite 200<br>Austin, Texas 78746<br>TBPE No. 1298 |  |

WEST

A



EAST

A'

EXPLANATION

MW-54 → WELL IDENTIFICATION

→ WELL

→ SCREEN INTERVAL

→ LITHOLOGIC CONTACTS

---▲--- POTENTIOMETRIC SURFACE MEASURED AUGUST-OCTOBER 2008

NOTE:  
SCREEN INTERVAL NOT AVAILABLE FOR MW-3, MW-40, MW-41, MW-44, RW-42



WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 10/24/11 FILE: WestRef-B115

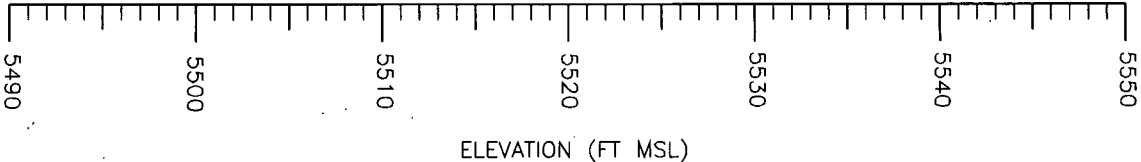
FIGURE 7

CROSS SECTION A-A'  
WEST TO EAST  
BLOOMFIELD REFINERY



Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298

# SOUTH 'B'

EXPLANATION

MMV-8 — WELL IDENTIFICATION

**F**

- SCREEN INTERVAL

LITHOLOGIC CONTACTS

SCALE IN FEET  
VERTICAL EXAGGERATION = 20X

POTENTIOMETRIC SURFACE MEASURED AUGUST-OCTOBER 2008

NOTE:  
SCREEN INTERVAL NOT AVAILABLE FOR MW-1, MW-3, MW-5



Legend

- Monitoring Well
- Observation Well
- Recovery Well
- Collection Well
- Piezometer
- Seep
- Site
- Approximate Property Line
- Groundwater Elevation Contours
- Inferred Groundwater Elevation
- Groundwater Flow Direction - Dashed where inferred
- Well ID  
Groundwater Elevation (ft amsl)

Notes:  
• Deeper Well; data not used to contour.

3rd Quarter  
August 17th

**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining | DATE: 10/24/11 | FILE: WestRef-B76

FIGURE 9

POTENTIOMETRIC SURFACE MAP  
AUGUST 2009  
BLOOMFIELD REFINERY

**RPS**  
Cleto Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TRPE No. 1298

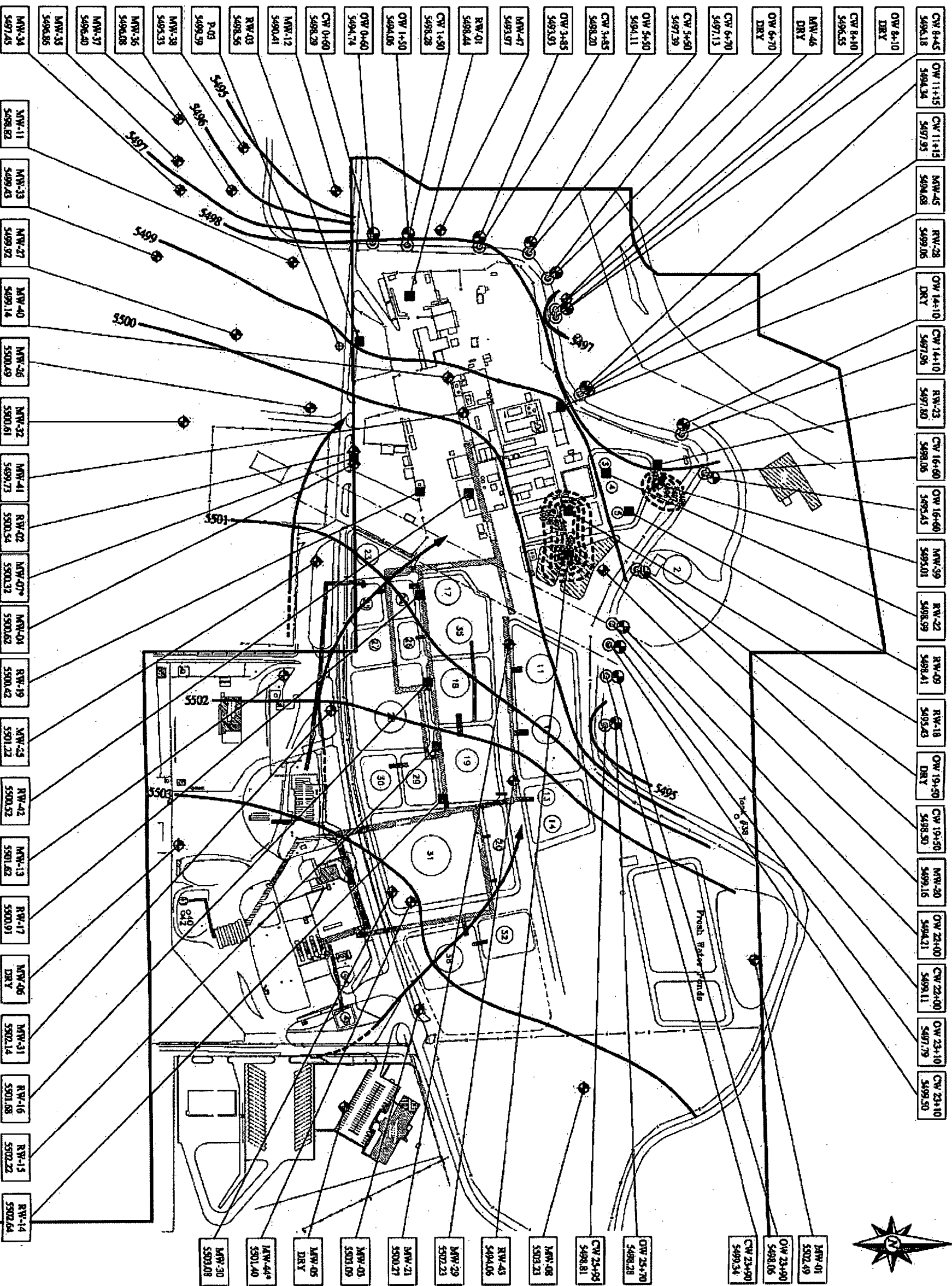
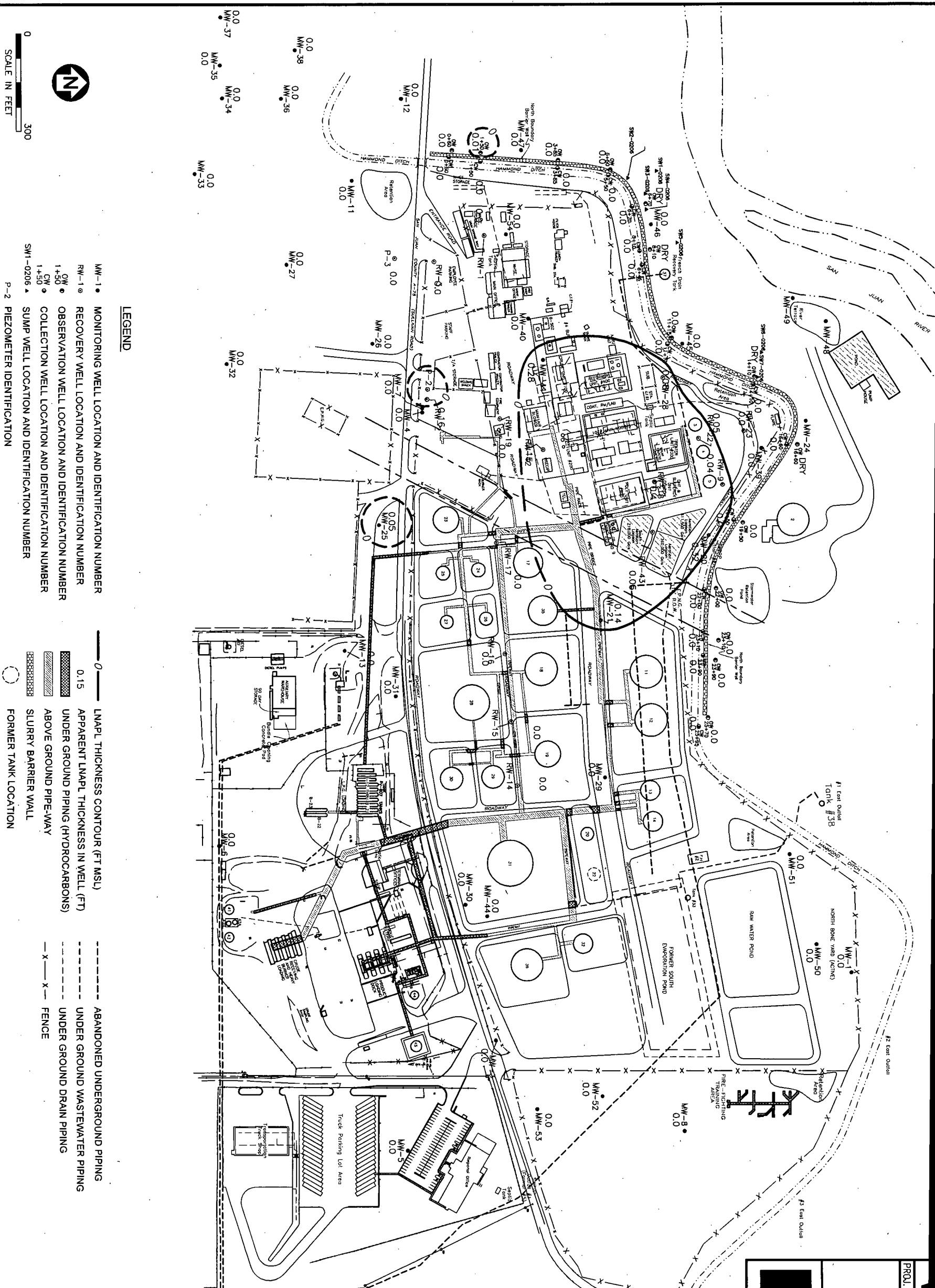


FIGURE 10  
SEPARATE PHASE  
HYDROCARBON THICKNESS MAP  
AUGUST 2008  
BLOOMFIELD REFINERY

**RPS**  
Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBE No. 1298



LEGEND

- MW-1 • MONITORING WELL LOCATION AND IDENTIFICATION NUMBER
- RW-1 • RECOVERY WELL LOCATION AND IDENTIFICATION NUMBER
- OW • OBSERVATION WELL LOCATION AND IDENTIFICATION NUMBER
- 1+50 • COLLECTION WELL LOCATION AND IDENTIFICATION NUMBER
- SM-0206 • SUMP WELL LOCATION AND IDENTIFICATION NUMBER
- P-2 • PIEZOMETER IDENTIFICATION
- 0 — LNAPL THICKNESS CONTOUR (FT MSU)
- 0.15 — APPARENT LNAPL THICKNESS IN WELL (FT)
- UNDER GROUND PIPING (HYDROCARBONS)
- ABOVE GROUND PIPE-WAY
- SLURRY BARRIER WALL
- FORMER TANK LOCATION
- ABANDONED UNDERGROUND PIPING
- UNDER GROUND WASTEWATER PIPING
- UNDER GROUND DRAIN PIPING
- FENCE



0 360  
APPROXIMATE SCALE IN FEET

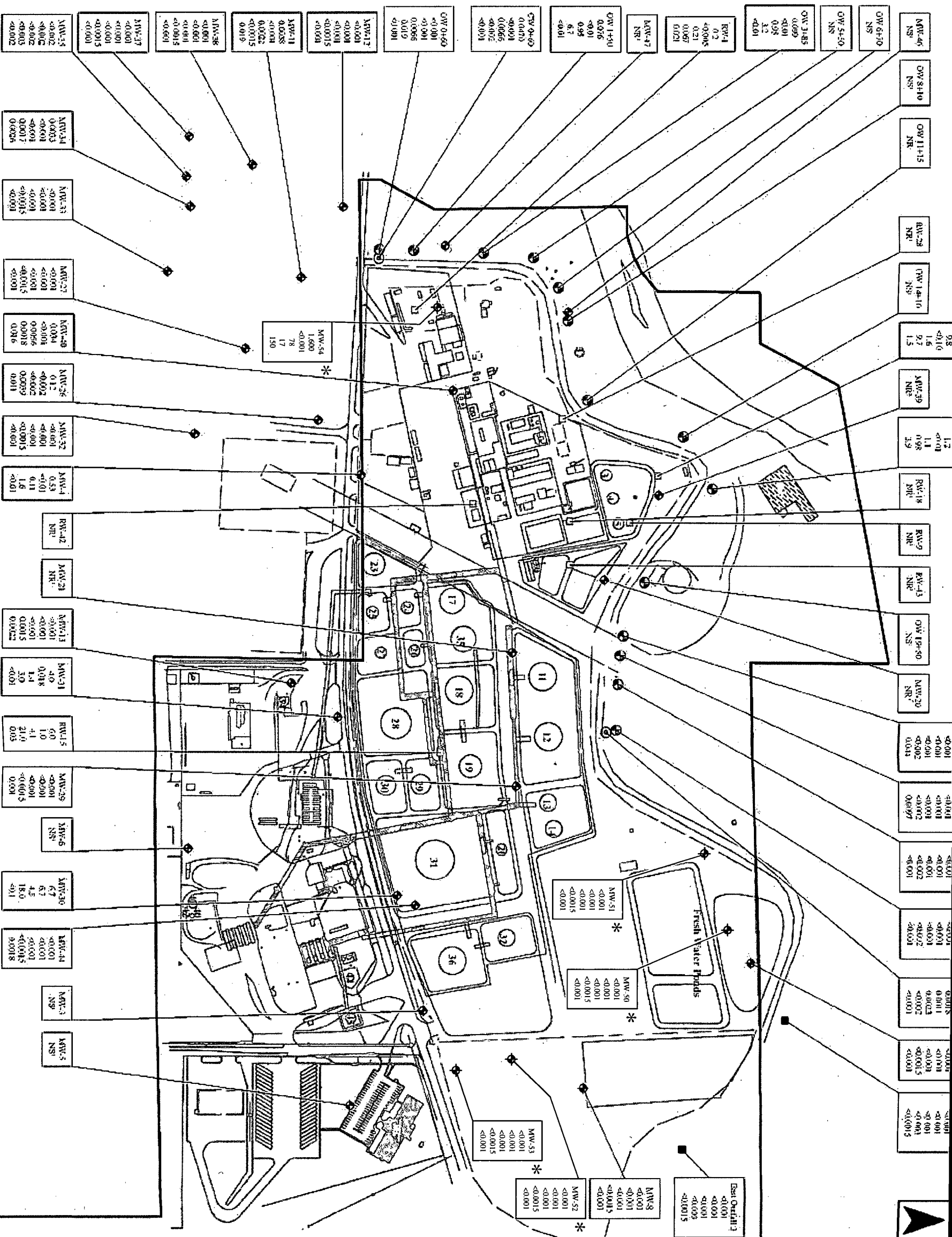
### Legend

- Monitoring Well
- Observation Well
- Recovery Well
- Collection Well
- Outfall
- Site
- Approximate Property Line

MW-38  
< 0.001  
- Well Id  
- Benzene  
< 0.001  
- Toluene  
- Ethylbenzene  
< 0.001  
- Xylenes, Total  
0.001  
- MTBE

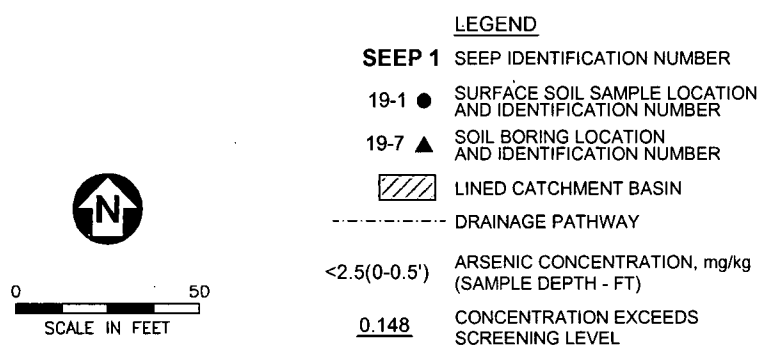
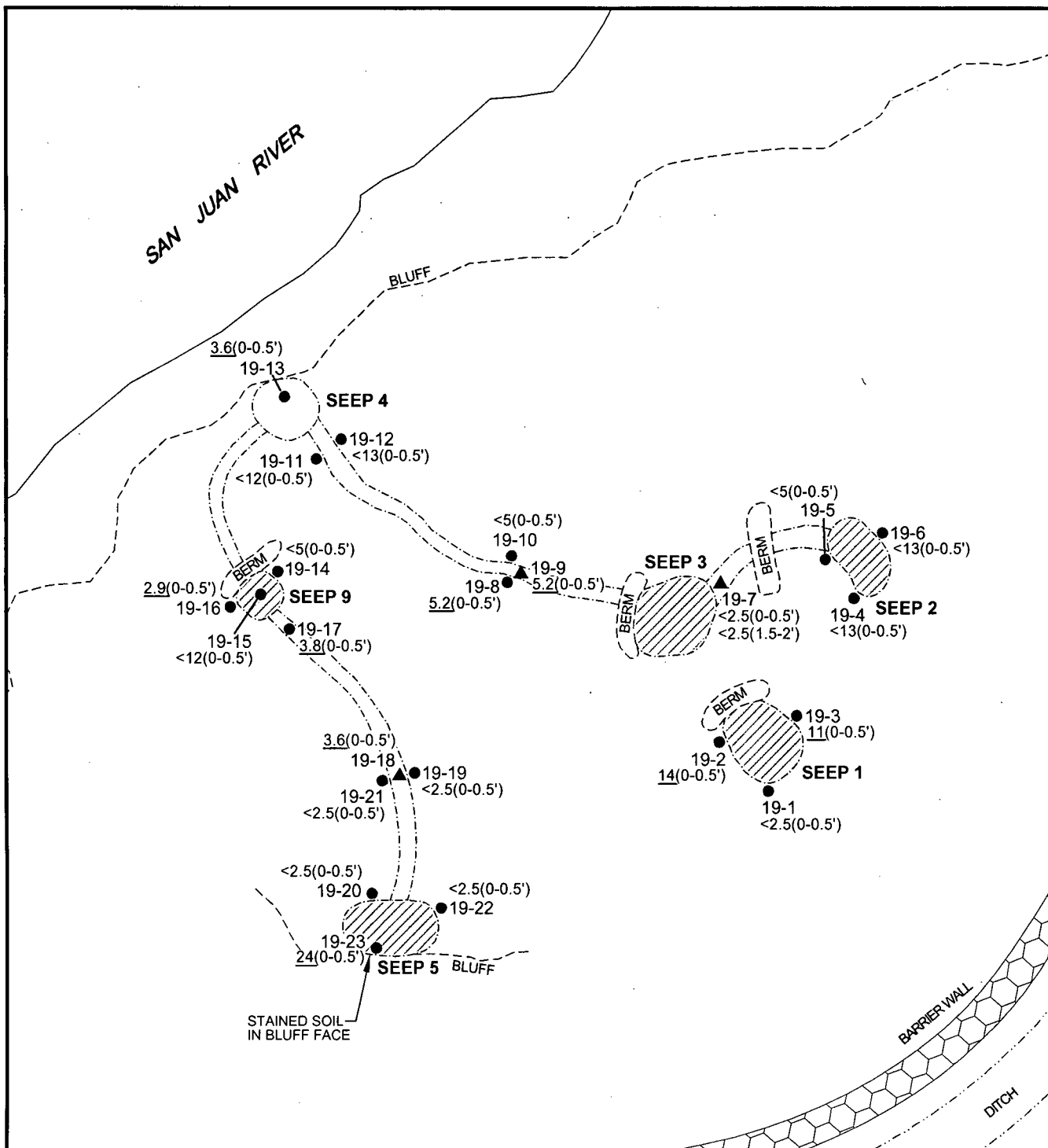
### Notes:

- All concentrations in milligrams per liter (mg/L)
- NS<sup>1</sup> = Well is Dry or Not Enough Water to Sample- No sample
- NS<sup>2</sup> = Sample Inadvertently not Collected this Sampling Event
- NR<sup>1</sup> = No Sample Required - Well Contains Separate Phase Hydrocarbon
- NR<sup>2</sup> = No Sample Required per OCD and NMED Conditions
- \* = October 2008, All Other Samples Collected August 2008



**Western Refining**  
WESTERN REFINING SOUTHWEST  
PROJ. NO.: Western Refining DATE: 10/24/11 FILE: WestRef-81118

**FIGURE 11**  
DISSOLVED-PHASE  
GROUNDWATER DATA  
AUGUST AND OCTOBER 2008  
BLOOMFIELD REFINERY  
Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
RPS  
TPE No. 1298



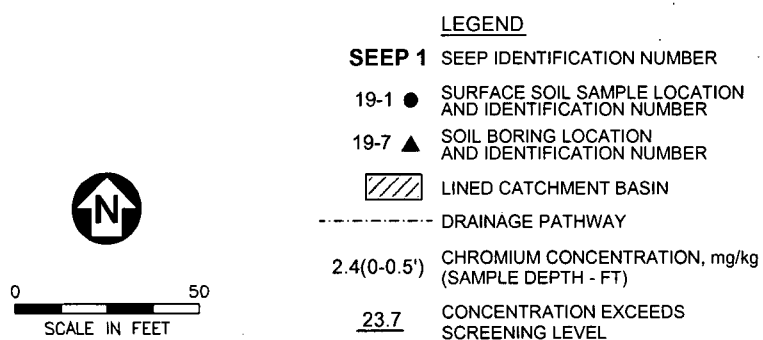
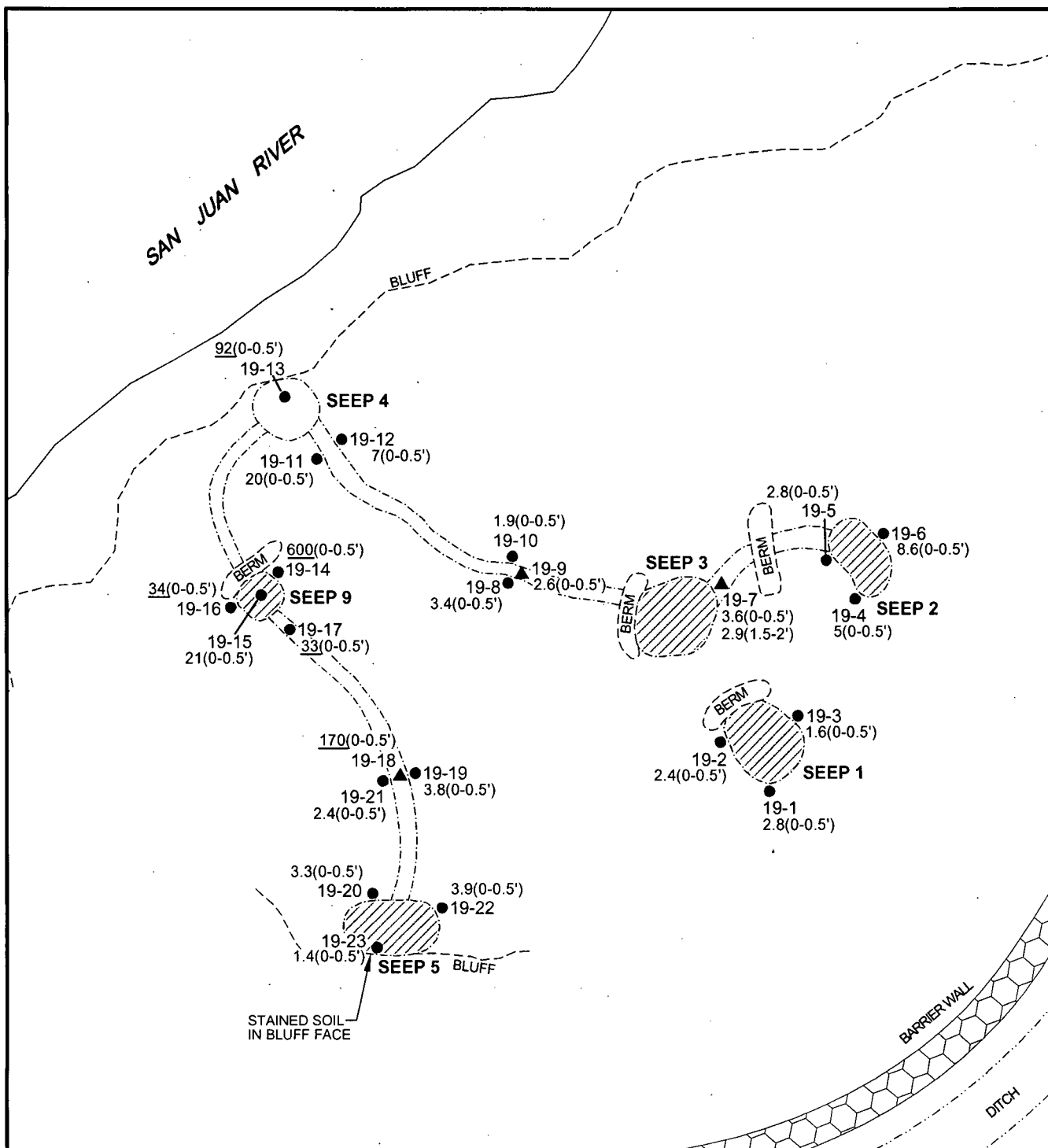
**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 10/24/11 FILE: WestRef-A136

**FIGURE 12**  
**AOC No. 19**  
**ARSENIC SOIL MAP**  
**BLOOMFIELD REFINERY**

**RPS**

Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298



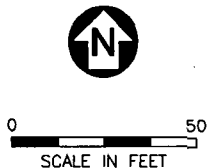
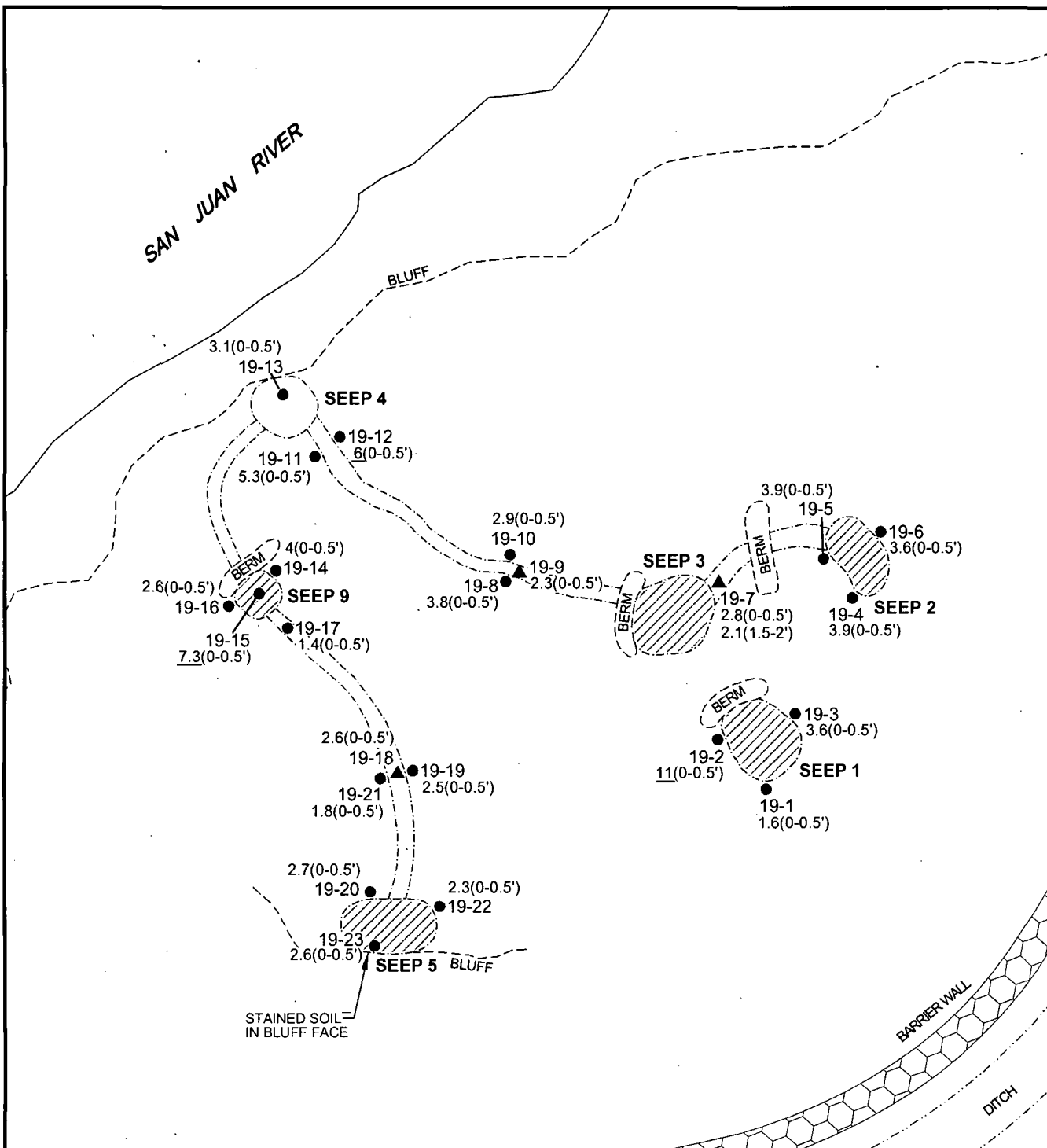
**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 10/24/11 FILE: WestRef-A136

**FIGURE 13**  
**AOC No. 19**  
**CHROMIUM SOIL MAP**  
**BLOOMFIELD REFINERY**

**RPS**

Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298



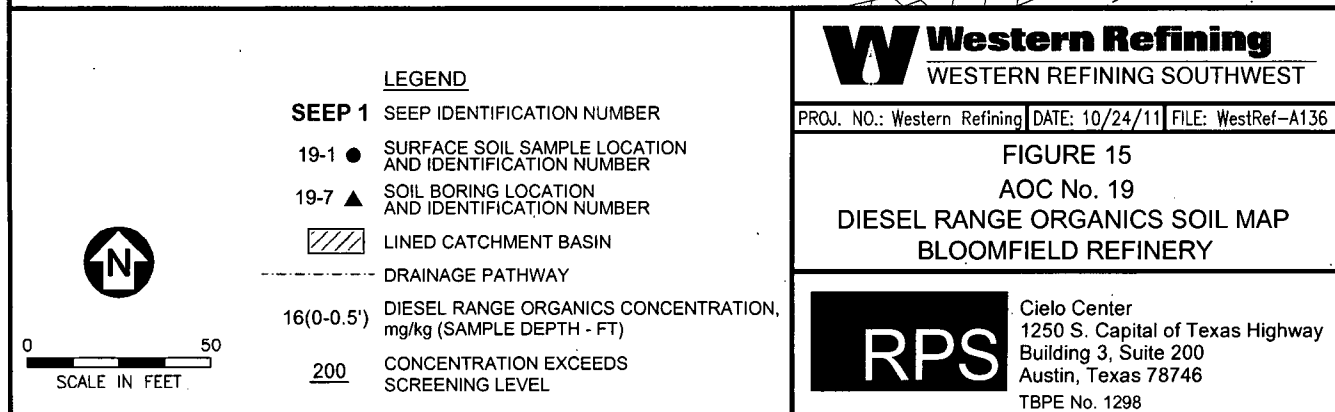
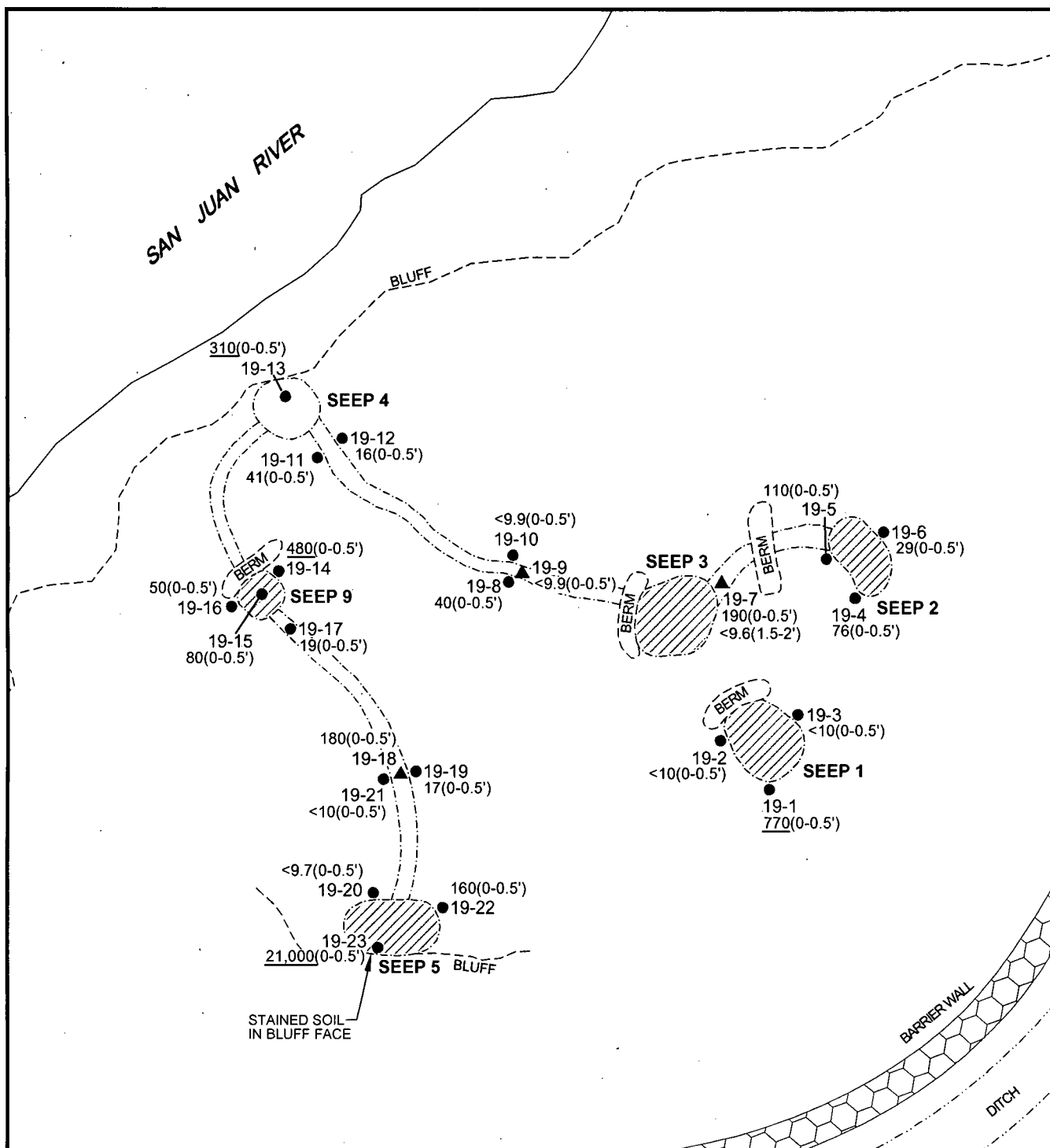
**Western Refining**  
WESTERN REFINING SOUTHWEST

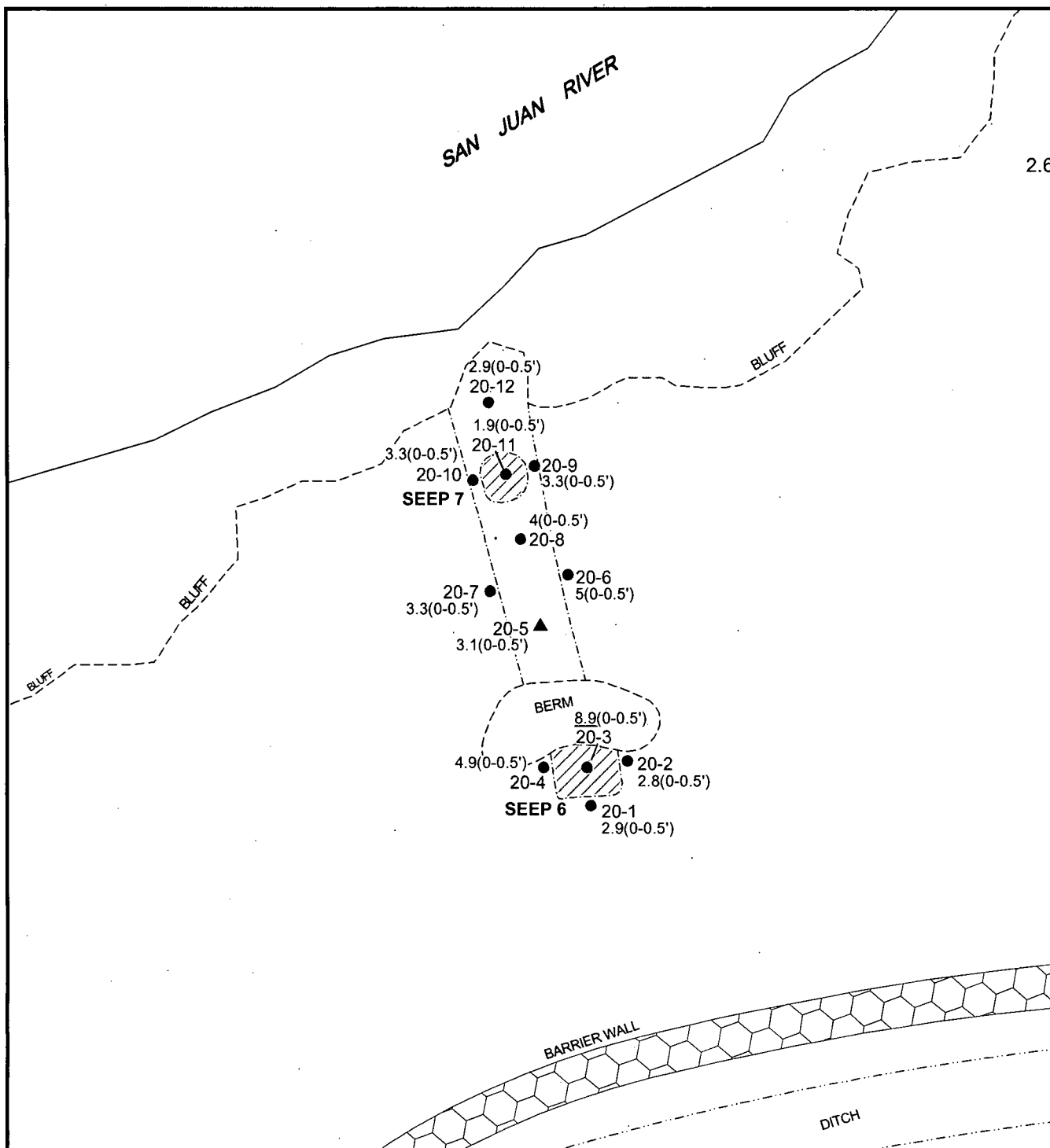
PROJ. NO.: Western Refining DATE: 10/24/11 FILE: WestRef-A136

**FIGURE 14**  
**AOC No. 19**  
**COBALT SOIL MAP**  
**BLOOMFIELD REFINERY**

**RPS**

Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298





#### LEGEND

**SEEP 6** SEEP IDENTIFICATION NUMBER

**20-1** ● SURFACE SOIL SAMPLE LOCATION AND IDENTIFICATION NUMBER

**20-5** ▲ SOIL BORING LOCATION AND IDENTIFICATION NUMBER

▨ LINED CATCHMENT BASIN

--- DRAINAGE PATHWAY

4.9(0-0.5') COBALT CONCENTRATION, mg/kg (SAMPLE DEPTH - FT)

5.51 CONCENTRATION EXCEEDS SCREENING LEVEL



0 40  
SCALE IN FEET



**Western Refining**

WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 10/24/11 FILE: WestRef-A136

**FIGURE 16**

**AOC No. 20**

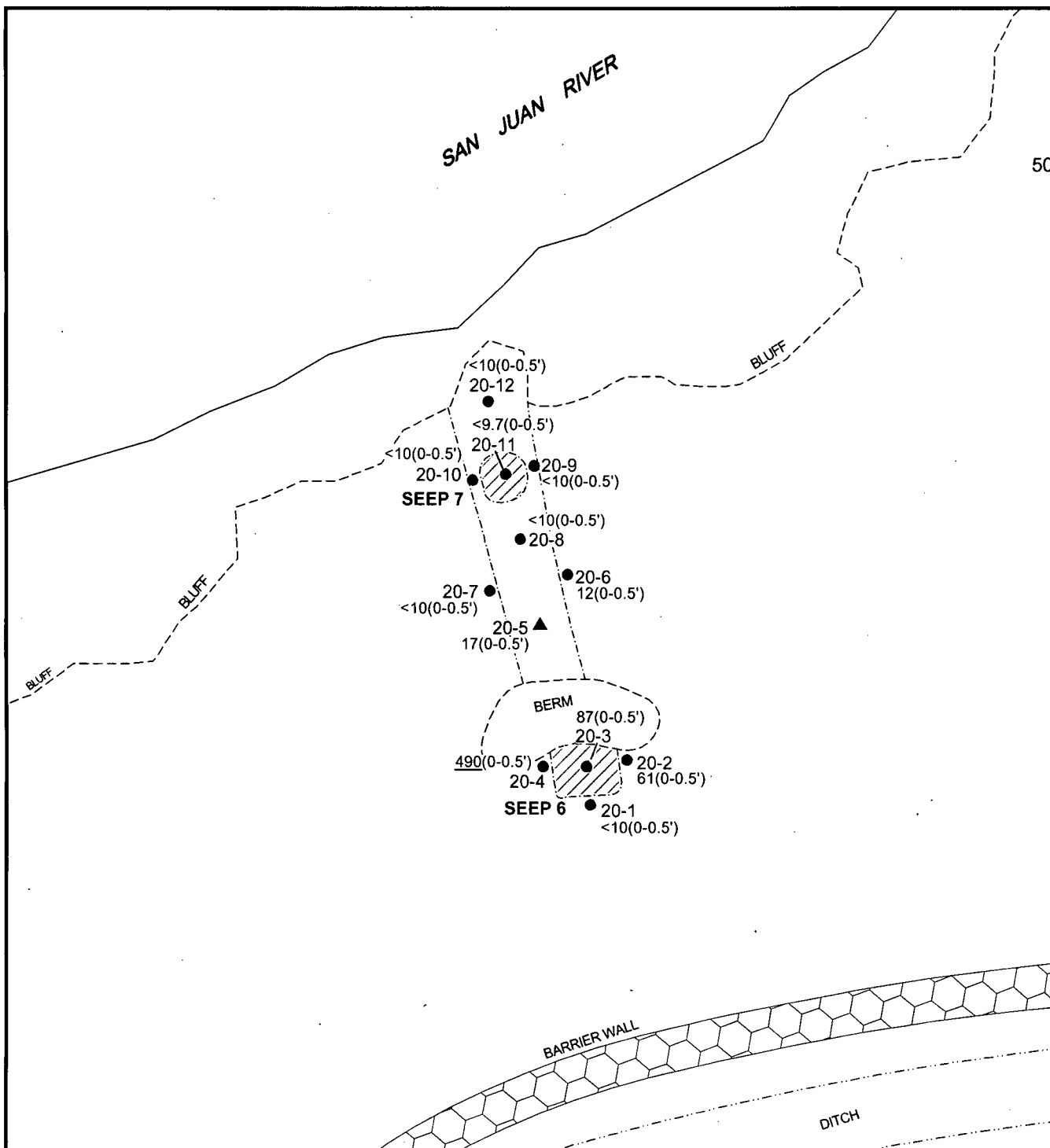
**COBALT SOIL MAP**

**BLOOMFIELD REFINERY**

**RPS**

Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298





#### LEGEND

**SEEP 6** SEEP IDENTIFICATION NUMBER

20-1 ● SURFACE SOIL SAMPLE LOCATION AND IDENTIFICATION NUMBER

20-5 ▲ SOIL BORING LOCATION AND IDENTIFICATION NUMBER

▨ LINED CATCHMENT BASIN

--- DRAINAGE PATHWAY

17(0-0.5') DIESEL RANGE ORGANICS CONCENTRATION, mg/kg (SAMPLE DEPTH - FT)

200 CONCENTRATION EXCEEDS SCREENING LEVEL



0 40  
SCALE IN FEET

**Western Refining**  
WESTERN REFINING SOUTHWEST

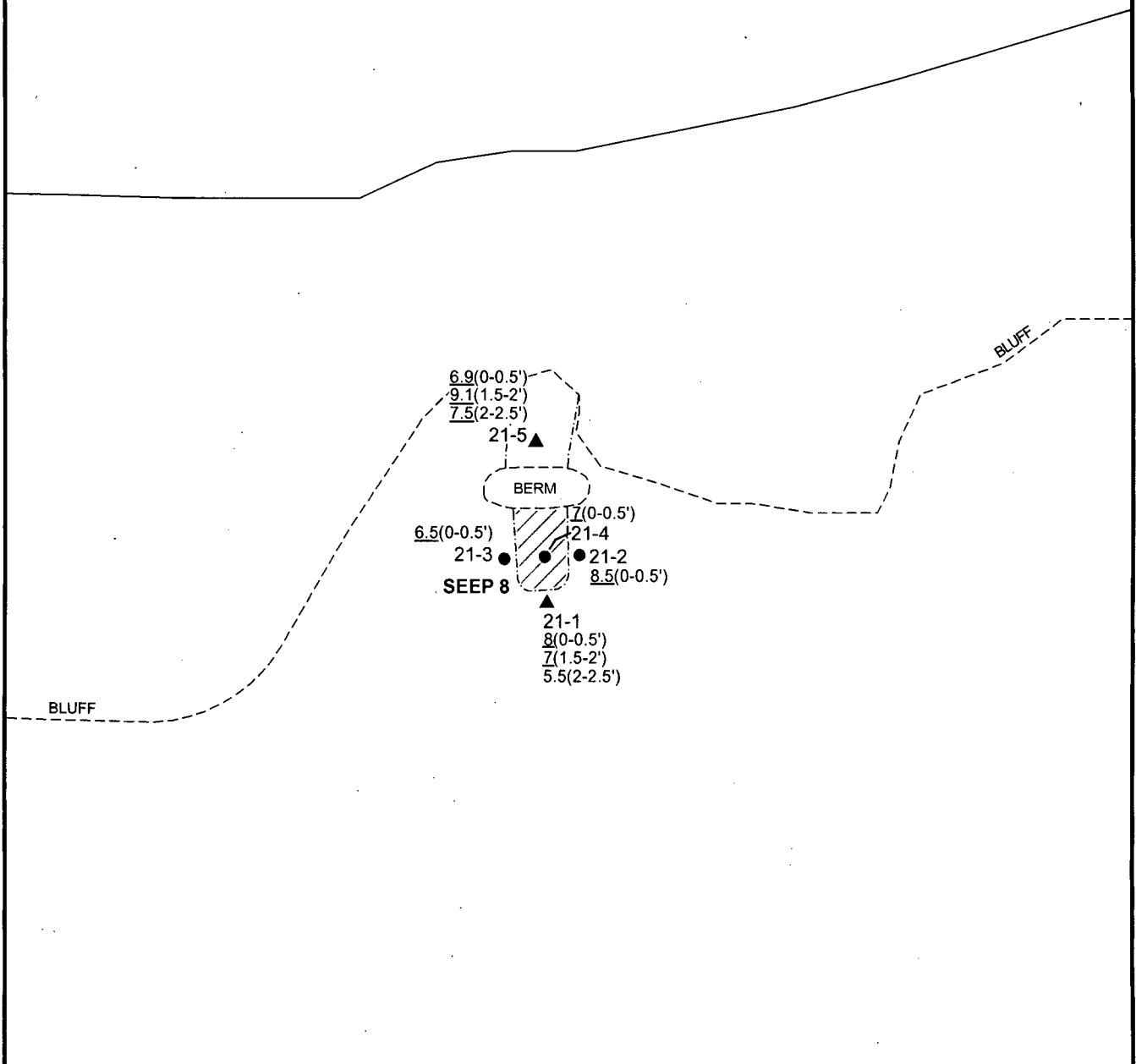
PROJ. NO.: Western Refining DATE: 10/24/11 FILE: WestRef-A136

**FIGURE 17**  
**AOC No. 20**  
**DIESEL RANGE ORGANICS SOIL MAP**  
**BLOOMFIELD REFINERY**



**RPS**

Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298

SAN JUAN RIVER



# LEGEND

- SEEP 8** SEEP IDENTIFICATION NUMBER
- 21-2 ● SURFACE SOIL SAMPLE LOCATION AND IDENTIFICATION NUMBER
- 21-1 ▲ SOIL BORING LOCATION AND IDENTIFICATION NUMBER
-  LINED CATCHMENT BASIN
-  DRAINAGE PATHWAY
- 5.5(0-0.5') COBALT CONCENTRATION, mg/kg (SAMPLE DEPTH - FT)
- 5.51 CONCENTRATION EXCEEDS SCREENING LEVEL



0 30  
SCALE IN FEET

**W Western Refining**  
WESTERN REFINING SOUTHWEST

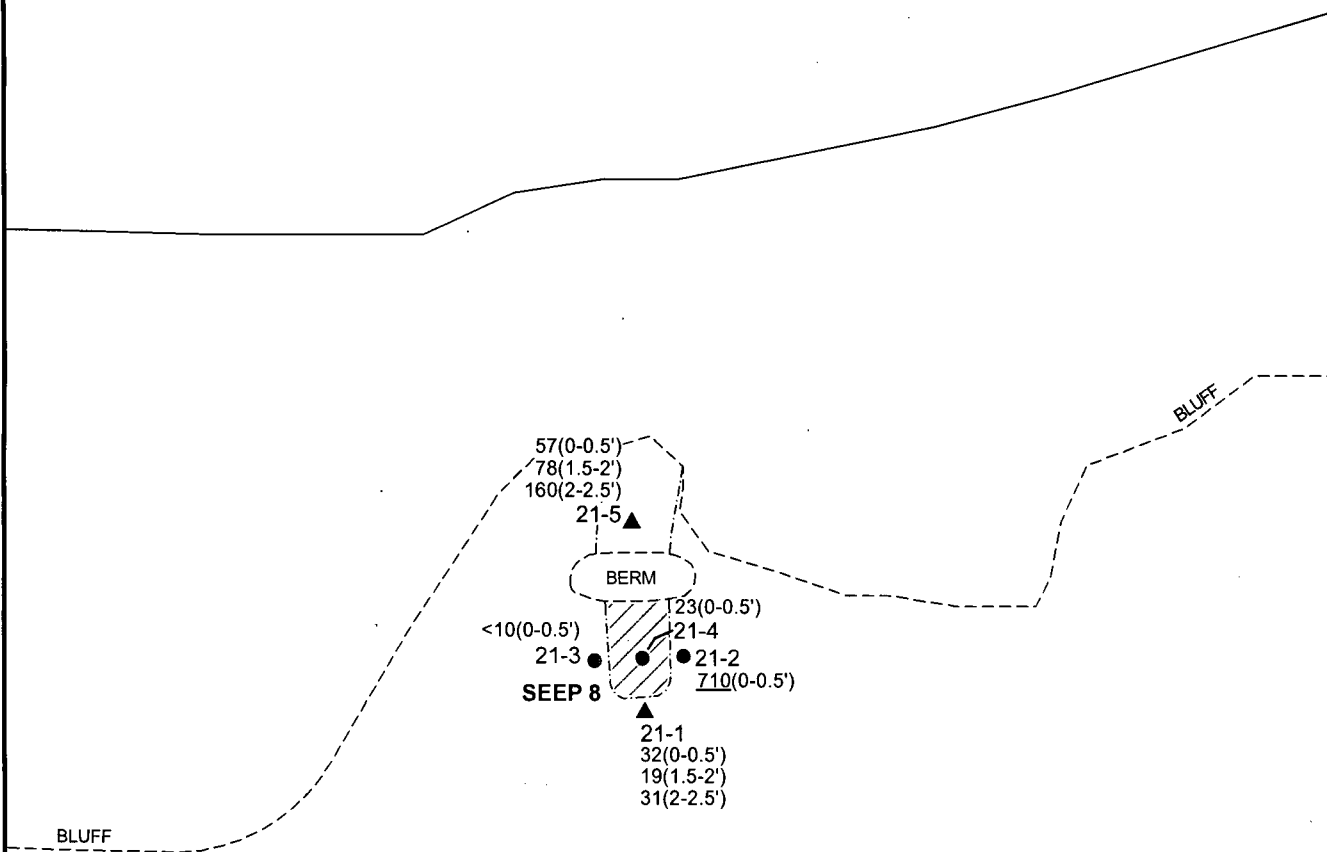
PROJ. NO.: Western Refining DATE: 10/24/11 FILE: WestRef-A136

FIGURE 18  
AOC No. 21  
COBALT SOIL MAP  
BLOOMFIELD REFINERY

**RPS**

Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298

SAN JUAN RIVER



**LEGEND**

**SEEP 8** SEEP IDENTIFICATION NUMBER

21-2 ● SURFACE SOIL SAMPLE LOCATION AND IDENTIFICATION NUMBER

21-1 ▲ SOIL BORING LOCATION AND IDENTIFICATION NUMBER

LINED CATCHMENT BASIN

DRAINAGE PATHWAY

32(0-0.5') DIESEL RANGE ORGANICS CONCENTRATION, mg/kg (SAMPLE DEPTH - FT)

200 CONCENTRATION EXCEEDS SCREENING LEVEL



0 30  
SCALE IN FEET

**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 10/24/11 FILE: WestRef-A136

**FIGURE 19**  
**AOC No. 21**  
**DIESEL RANGE ORGANICS SOIL MAP**  
**BLOOMFIELD REFINERY**

**RPS**

Cielo Center  
1250 S. Capital of Texas Highway  
Building 3, Suite 200  
Austin, Texas 78746  
TBPE No. 1298

# **Appendix A**

---

## **Field Methods**

## **Field Methods**

Pursuant to Section IV of the Order, an investigation of soils was conducted to determine and evaluate the presence, nature, extent, fate, and transport of contaminants. To accomplish this objective, soil was collected from borings and surface sample locations at AOCs No. 19, 20, and 21. The field methods are described below and individual discussions are presented for the following activities:

- Drilling procedures;
- Soil screening;
- Decontamination procedures;
- Sample collection and handling procedures;
- Equipment calibration; and
- Management of investigation derived waste.

### **Drilling Procedures**

The soil borings were completed using a hand auger. Soil samples were collected continuously and logged by a qualified geologist in accordance with the Unified Soil Classification System (USCS) nomenclature. As shown on the boring logs, the data recorded included the lithologic interval, symbol, percent recovery, field screening results, and a sample description of the cuttings and core samples.

### **Soil Screening**

Samples obtained from the borings and surface samples were screened in the field on 2-foot intervals for evidence of contaminants. Field screening results were recorded on the soil boring logs. Field screening results were used to aid in the selection of soil samples for laboratory analysis. The primary screening methods include: (1) visual examination, (2) olfactory examination, and (3) headspace vapor screening for volatile organic compounds.

Visual screening included examining the soil samples for evidence of staining caused by petroleum-related compounds or other substances that may have caused staining of soils such as elemental sulfur or cyanide compounds. Headspace vapor screening was conducted and involved placing a soil sample in a plastic sealable bag allowing space for ambient air. The bag was sealed, labeled and then shaken gently to expose the soil to the air trapped in the container. The sealed bag was allowed to rest for a minimum of 5 minutes while the vapors equilibrated. Vapors present within the sample bag's headspace were then measured by inserting the probe of a MiniRae 3000 portable volatile organic constituent (VOC) monitor PGM-7600 in a small opening in the bag. The maximum value and the ambient air temperature were

recorded on the field boring log for each sample. The screening results are presented in Table 2. Field screening results and any conditions that were considered to be capable of influencing the results of the field screening were recorded on the field logs.

#### Decontamination Procedures

The sampling equipment coming in direct contact with the samples (e.g., hand augers and stainless steel trowel) were decontaminated using a brush, as necessary, to remove larger particulate matter followed by a rinse with potable water, wash with nonphosphate detergent, rinse with potable water, and double rinse with deionized water. In the event that more than one AOC was investigated during the day a new batch of wash water and rinse water was prepared prior to decontamination.

#### Sample Collection and Handling Procedures

Soil samples were collected using a stainless steel trowel or directly from the auger bucket for borings completed with a hand auger. The selected portion of the sample interval was placed in pre-cleaned, laboratory-prepared sample containers for laboratory chemical analysis. Three soil samples were collected for VOC analysis. An Encore® Sampler was used for collection of soil samples for low-level VOC analysis pursuant to EPA method 5035; the second sample aliquot (approximately 1 gram) was placed in a laboratory-prepared container with a methanol preservative; and the third sample aliquot was placed in an 8-ounce glass jar, which was filled to the top to minimize any head space.

Samples were immediately placed into an ice chest with ice. The samples were maintained in the custody of the sampler until the chain-of-custody form was completed and the ice chest was sealed for shipment to the laboratory.

#### Equipment Calibration

Soil vapor screening was conducted using a MiniRae 3000 portable VOC monitor PGM-7600. The instrument was calibrated at the beginning of each work day to a concentration of 100 ppm isobutylene.

#### Management of Investigation Derived Waste

The decontamination water from the drilling equipment was collected in five-gallon buckets and was subsequently placed in open top 55-gallon drums, which were sealed at the end of each work day. The decontamination water generated from sampling equipment was collected in

buckets and placed in open top 55-gallon drums, which were sealed at the end of each work day. The decon water was disposed in the Refinery's wastewater treatment system up-stream of the API Separator. Soil cuttings were also placed into an open-top 55-gallon drum, which was sealed when not in use. The drum of soils was labeled and temporarily stored in a concrete curbed area pending waste characterization and disposal.

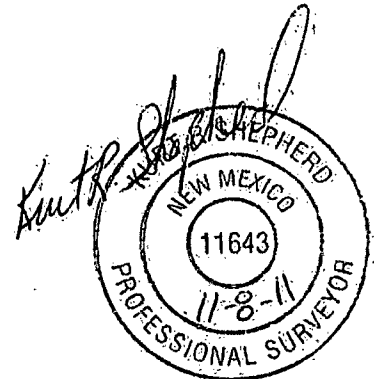


**Western Refining Southwest  
Seep Areas of Concern Field Ties**

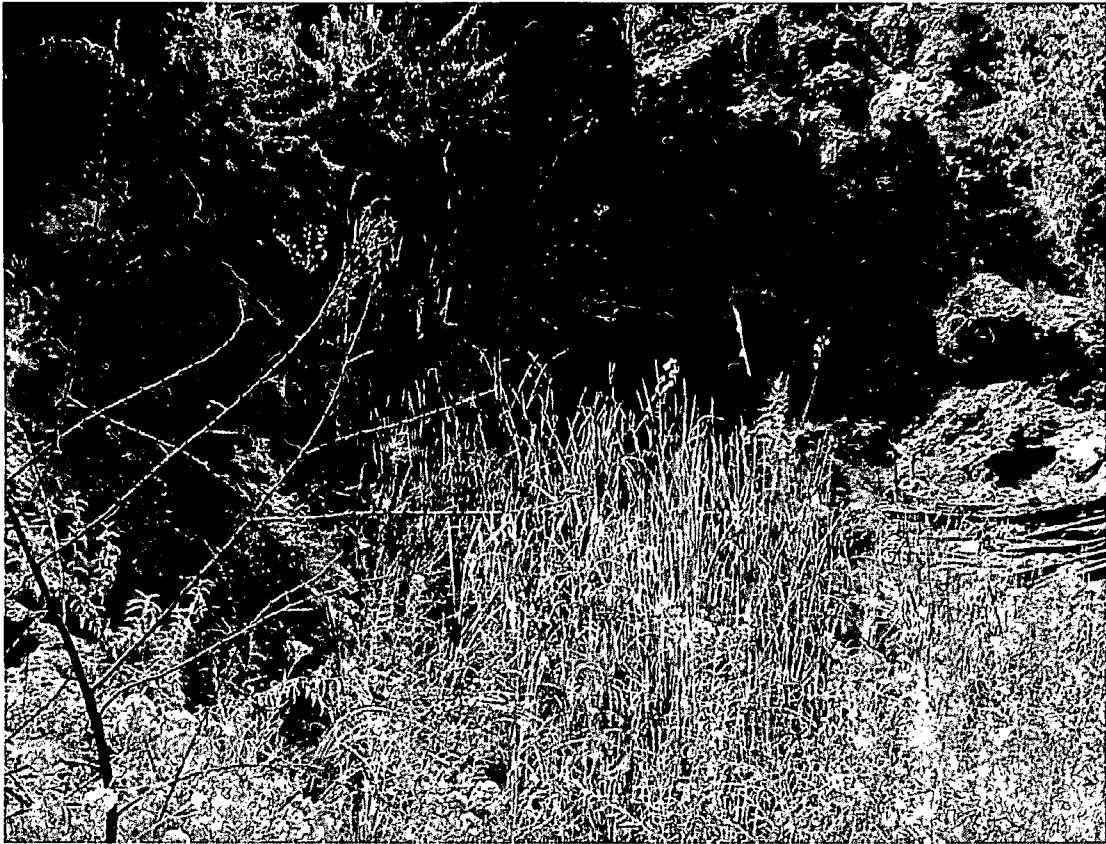
10/27/2011

Note\*\* This X, Y, Z shown  
here is derived from the  
City of Bloomfield  
Control Network

| Point | Northing  | Easting    | Elevation | Location  |
|-------|-----------|------------|-----------|-----------|
| 5000  | 74568.654 | 181637.006 | 5474.453  | AOC 19-1  |
| 5001  | 74591.068 | 181633.472 | 5468.928  | AOC 19-3  |
| 5002  | 74582.146 | 181619.347 | 5470.37   | AOC 19-2  |
| 5003  | 74648.195 | 181652.866 | 5470.392  | AOC 19-5  |
| 5004  | 74648.531 | 181659.695 | 5471.97   | AOC 19-4  |
| 5005  | 74660.661 | 181659.367 | 5472.586  | AOC 19-6  |
| 5006  | 74643.39  | 181609.048 | 5459.017  | AOC 19-7  |
| 5007  | 74641.537 | 181514.208 | 5443.684  | AOC 19-8  |
| 5008  | 74643.172 | 181515.364 | 5443.599  | AOC 19-9  |
| 5009  | 74644.548 | 181517.072 | 5443.81   | AOC 19-10 |
| 5010  | 74680.844 | 181466.492 | 5439.073  | AOC 19-11 |
| 5011  | 74687.795 | 181474.856 | 5439.119  | AOC 19-12 |
| 5012  | 74702.828 | 181462.667 | 5437.428  | AOC 19-13 |
| 5013  | 74629.846 | 181451.326 | 5445.7    | AOC 19-15 |
| 5014  | 74628.77  | 181446.148 | 5447.474  | AOC 19-16 |
| 5015  | 74622.64  | 181451.735 | 5446.746  | AOC 19-17 |
| 5016  | 74631.436 | 181457.65  | 5448.15   | AOC 19-14 |
| 5017  | 74574.618 | 181492.695 | 5461.77   | AOC 19-21 |
| 5018  | 74575.084 | 181497.765 | 5460.765  | AOC 19-18 |
| 5019  | 74575.963 | 181503.283 | 5461.075  | AOC 19-19 |
| 5020  | 74532.91  | 181482.057 | 5477.244  | AOC 19-20 |
| 5021  | 74530.609 | 181493.6   | 5477.52   | AOC 19-22 |
| 5022  | 74513.176 | 181482.442 | 5480.975  | AOC 19-23 |
| 5023  | 74472.531 | 181289.678 | 5467.509  | AOC 20-2  |
| 5024  | 74470.728 | 181275.869 | 5467.38   | AOC 20-4  |
| 5025  | 74471.467 | 181280.71  | 5465.067  | AOC 20-3  |
| 5026  | 74461.182 | 181283.544 | 5469.417  | AOC 20-1  |
| 5027  | 74513.196 | 181276.971 | 5456.308  | AOC 20-5  |
| 5028  | 74526.439 | 181278.604 | 5453.998  | AOC 20-6  |
| 5029  | 74530.099 | 181266.994 | 5453.359  | AOC 20-7  |
| 5030  | 74535.077 | 181273.44  | 5451.499  | AOC 20-8  |
| 5031  | 74540.095 | 181268.539 | 5450.617  | AOC 20-10 |
| 5032  | 74540.076 | 181274.717 | 5449.867  | AOC 20-11 |
| 5033  | 74537.171 | 181282.135 | 5451.334  | AOC 20-9  |
| 5034  | 74555.04  | 181272.202 | 5446.088  | AOC 20-12 |
| 5035  | 74450.302 | 181046.206 | 5473.066  | AOC 21-1  |
| 5036  | 74460.204 | 181050.86  | 5472.78   | AOC 21-2  |
| 5037  | 74472.412 | 181047.834 | 5470.482  | AOC 21-5  |
| 5038  | 74460.782 | 181045.136 | 5471.458  | AOC 21-4  |
| 5039  | 74465.222 | 181040.544 | 5471.977  | AOC 21-3  |







AOC No. 19 – Seep and Catchment Basin No. 1



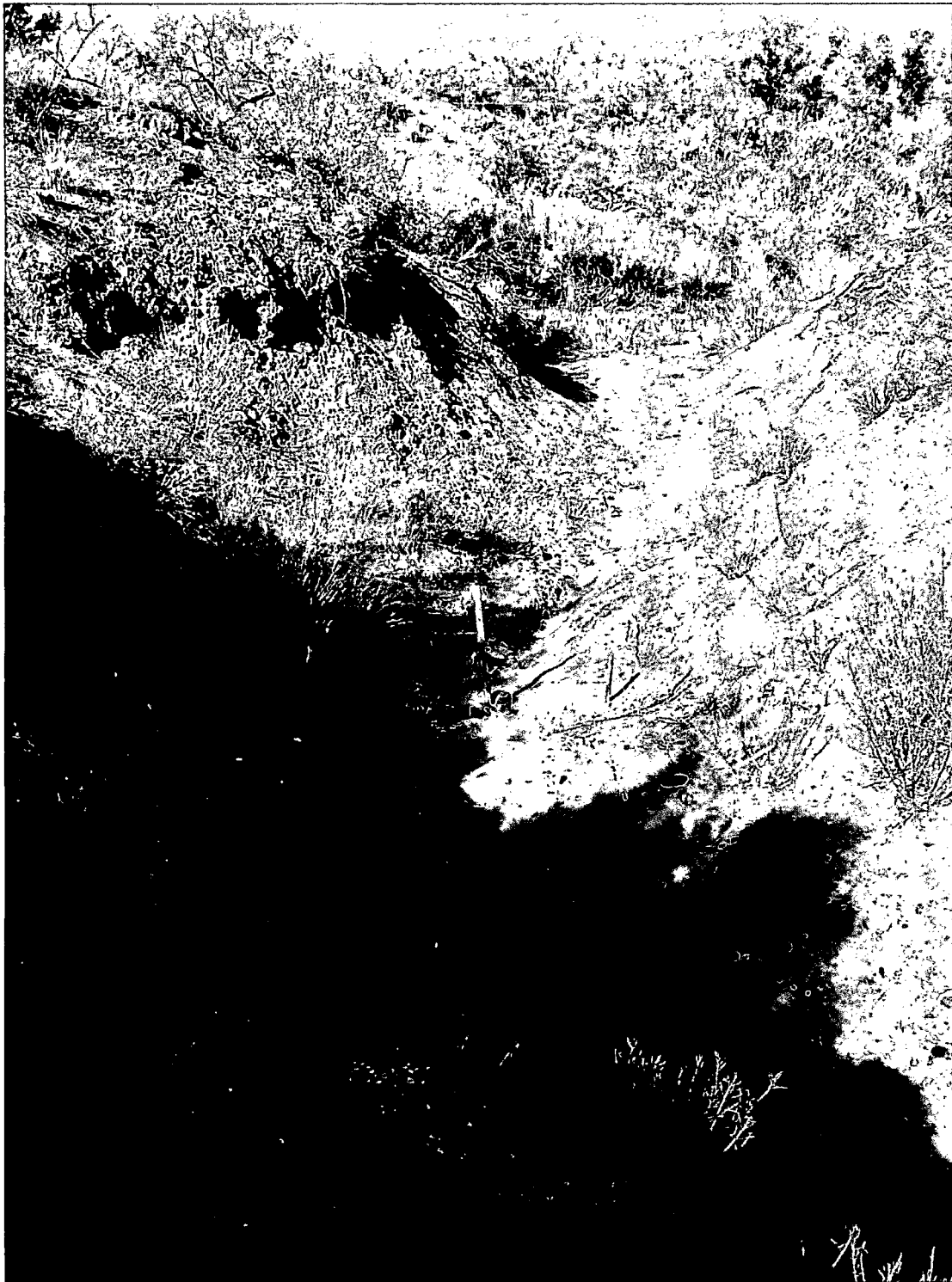
AOC No. 19 – Seep No. 2



AOC No. 19 – Sediment Within Catchment Basin No. 3



AOC No. 19 – Seep 9 Looking Southwest



AOC No. 19 – Looking Northwest Toward Catchment No. 4 from Catchment No. 3



AOC No. 19 – Looking northward at Seep 4



AOC No. 19 – Between Seep 5 and Seep 9





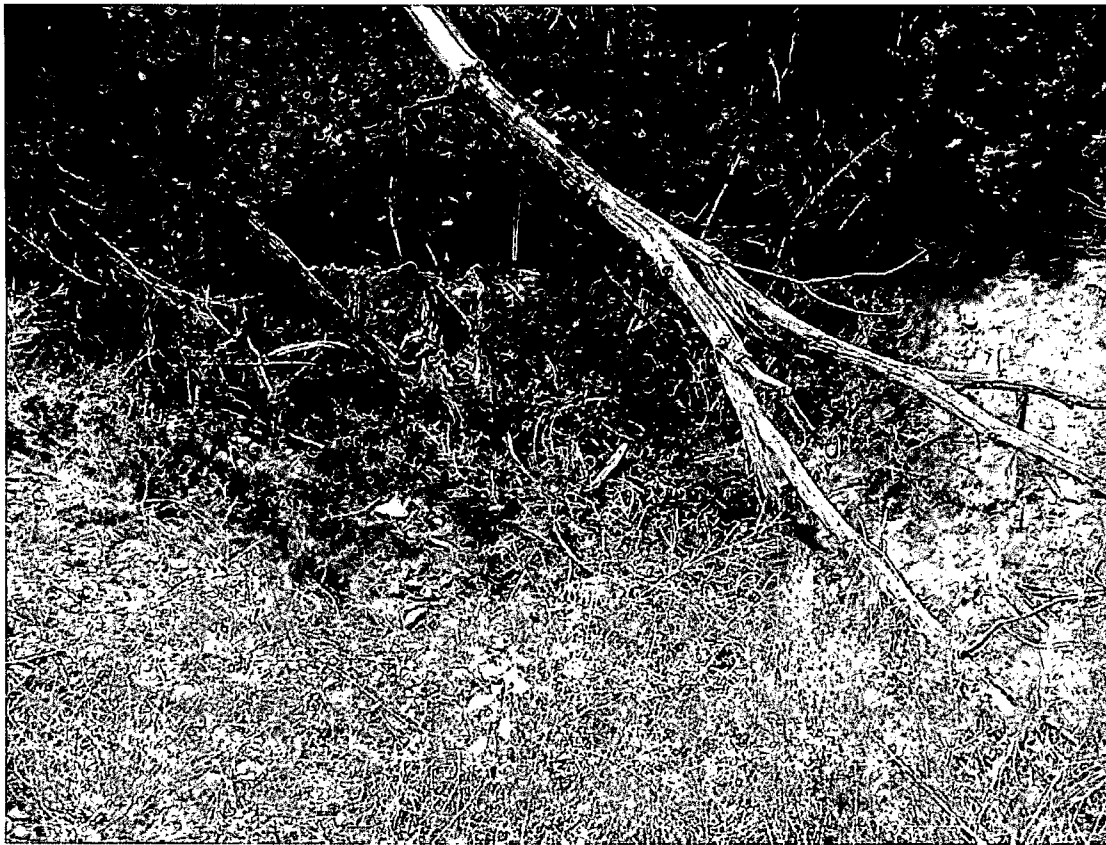
AOC No. 19 – HDPE Liner of Seep No. 5 Catchment Basin



AOC No. 19 – Stained Soil in Bluff Face of Seep No. 5



AOC No. 19 – Location of sample at Seep No. 5.



AOC 20 – Seep No. 6 With Water in Catchment Basin No. 6



AOC 20 – Looking North Toward Seep No. 7 From Berm Below Catchment Basin No. 6





AOC 20 – Seep No. 7 and Bluff Overlooking San Juan River



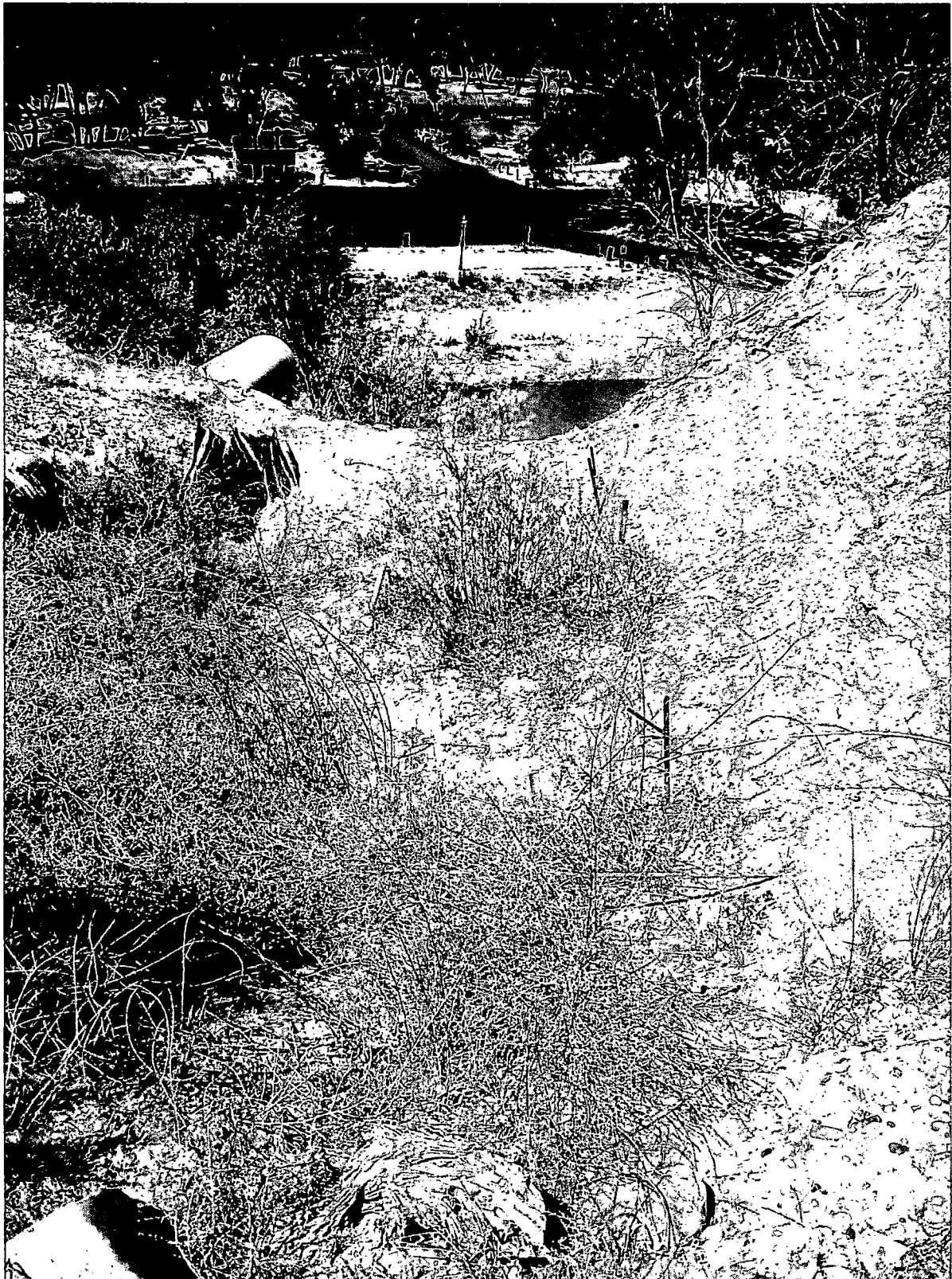
AOC No. 20 – From Seep No. 7, looking south.



AOC No. 20 – Between Seep No. 7 and Seep No. 6, looking south.



AOC 21 – Seep No. 8 (Below Wood Wall) and Liner of Catchment Basin No. 8



AOC No. 21 – Looking north from above Seep No. 8.



AOC 21 – Looking North Over San Juan River Bluff From Catchment Basin No. 8

Google maps

bloomfield, NM

Find businesses, addresses and places of interest.

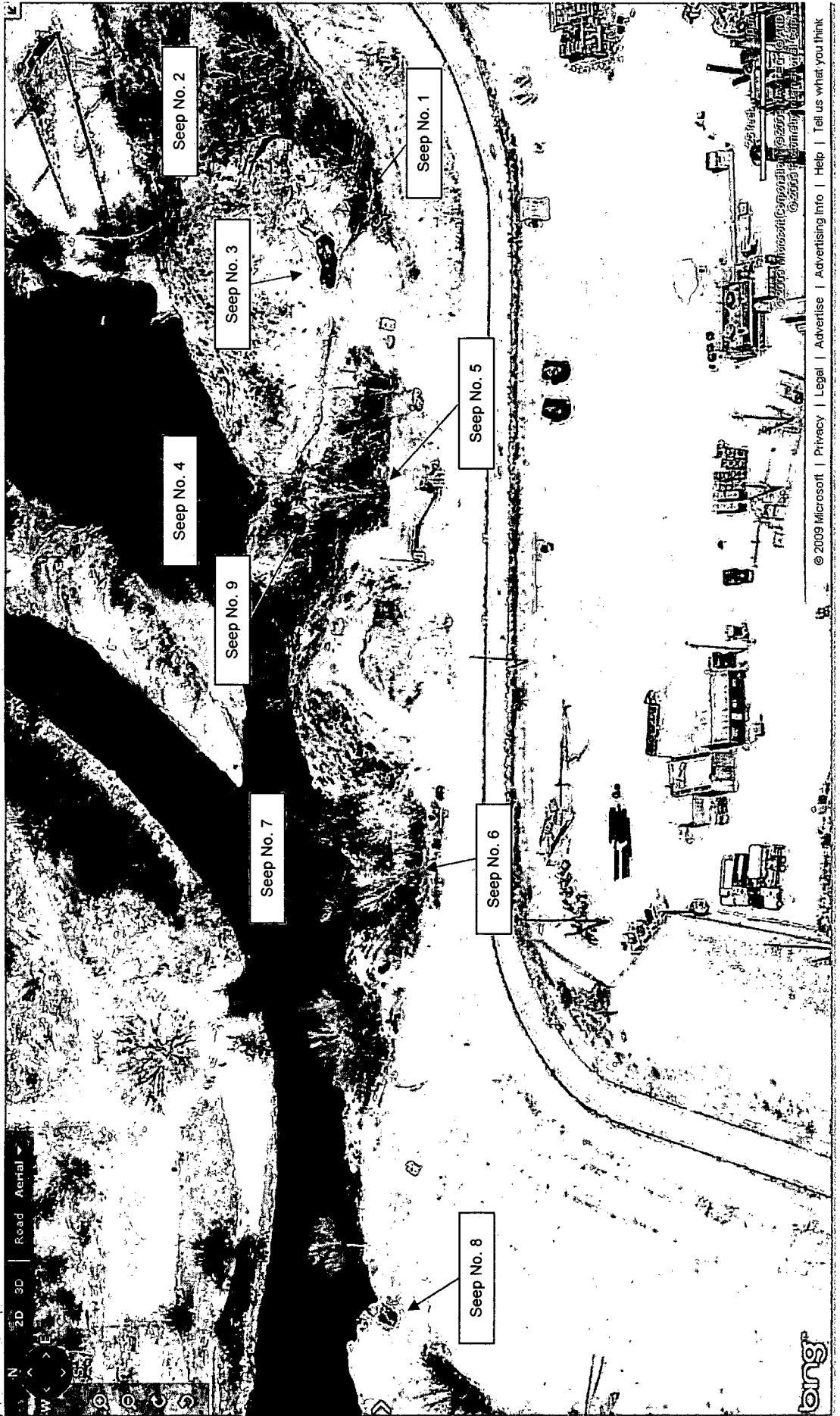
Search Maps

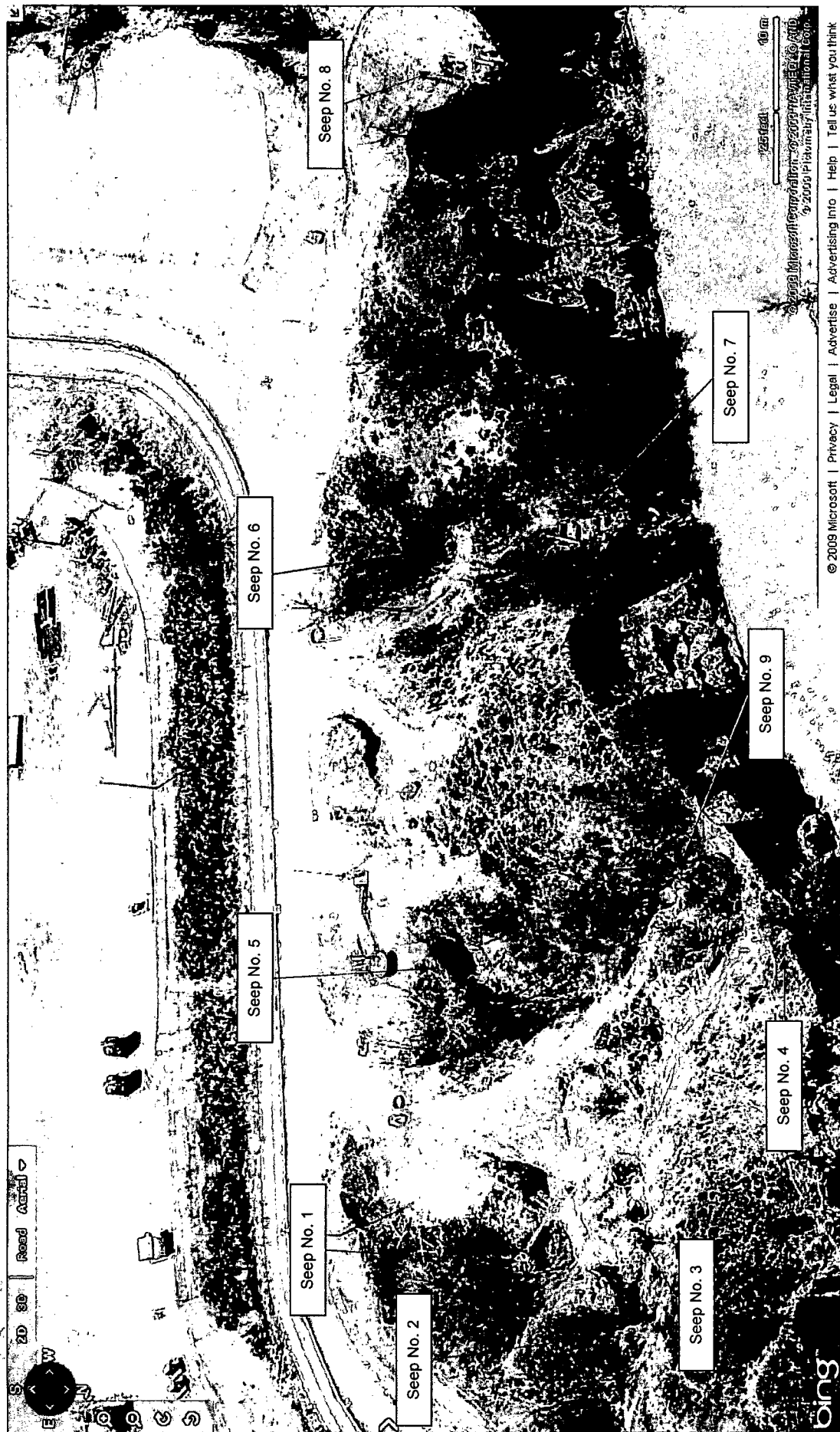
Show search options



© 2006 Google. All rights reserved. Terms of Service













**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.957' W 107°58.667'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5474.453  
**Site Coordinates:**  
**N** 74568.654      **E** 181637.006

**Boring No.:** AOC 19-1  
**Start Date:** 8/15/2011 12:00  
**Finish Date:** 8/15/2011 12:30

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1220 | G/2V/<br>2E/3J                |            | 92<br>95°F             |            | 100          | Ground Surface<br><b>Clayey Sand (SC)</b><br>Fine grain, compact, moist to saturated, gray to black, 1/2 inch to +3 inch gravel present, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.958' W 107°58.670'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5470.37  
**Site Coordinates:**  
**N** 74582.146      **E** 181619.347

**Boring No.:** AOC 19-2  
**Start Date:** 8/15/2011 12:30  
**Finish Date:** 8/15/2011 13:00

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1240 | G/2V/<br>2E/3J                |            | 5.0<br>95°F            |            | 100          | Ground Surface<br><b>Clayey Sand (SC)</b><br>Fine grain, compact, dry to damp, tan, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.960' W 107°58.669'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5468.928  
**Site Coordinates:**  
**N** 74591.068      **E** 181633.472

**Boring No.:** AOC 19-3  
**Start Date:** 8/15/2011 11:50  
**Finish Date:** 8/15/2011 12:00

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1158 | G/2V/<br>2E/3J                |            | 5.0<br>95°F            |            | 100          | Ground Surface<br><b>Clayey Sand (SC)</b><br>Fine grain, compact, dry to damp, tan, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.971' W 107°58.663'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5471.97  
**Site Coordinates:**  
**N** 74648.531      **E** 181659.695

**Boring No.:** AOC 19-4  
**Start Date:** 8/15/2011 13:20  
**Finish Date:** 8/15/2011 13:40

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 1325 | G/2V/<br>2E/3J                |            | 4.2<br>99°F            |            | 100          | Ground Surface<br><b>Clayey Sand (SC)</b><br>Very fine grain, loose, dry, light brown, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.974' W 107°58.665'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5470.392  
**Site Coordinates:**  
**N** 74648.195      **E** 181652.866

**Boring No.:** AOC 19-5  
**Start Date:** 8/15/2011 13:40  
**Finish Date:** 8/15/2011 14:00

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1345 | G/2V/<br>2E/3J                |            | 3.5<br>99°F            |            | 100          | Ground Surface<br><b>Silty Sand (SM)</b><br>Very fine grain, loose, dry, light brown, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.971' W 107°58.665'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5472.586  
**Site Coordinates:**  
**N** 74660.661      **E** 181659.367

**Boring No.:** AOC 19-6  
**Start Date:** 8/15/2011 14:00  
**Finish Date:** 8/15/2011 14:10

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1405 | G/2V/<br>2E/3J                |            | 2.0<br>99°F            |            | 100          | Ground Surface<br><b>Clayey Sand (SC)</b><br>Very fine grain, loose, dry, light gray, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |



# LOG OF BORING

**Client:** Western Refining Southwest, Inc.

**Site:** SWMU Group #6, Bloomfield Refinery

**Job No.:** UEC01318 - Bloomfield, NM

**Geologist:** Tracy Payne

**Driller:** N/A

**Drilling Rig:** N/A

**Drilling Method:** N/A

**Sampling Method:** Stainless Steel Trowel

**Comments:** GPS Coordinates - N 36°41.969' W 107°58.672'

**Total Depth: 2' bgl**

**Ground Water:** Not Encountered

**Elev., TOC (ft. msl): --**

**Elev., PAD (ft. msl): --**

**Elev., GL (ft. msl): 5459.017**

**Site Coordinates:**

**N 74643.39**

**E 181609.048**

**Boring No.:** AOC 19-7

**Start Date:** 8/16/2011 12:55

**Finish Date:** 8/16/2011 13:15

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1300 | G/2V/<br>2E/3J                |            | 2.3<br>95°F            |            | 100          | Ground Surface  | 0           |
|             |              |      |                               |            | 2.0<br>95°F            |            | 100          | <b>Sand (SP)</b><br>Similar to above, no odor                                 |             |
|             | 1.5-2'       | 1305 | G/2V/<br>2E/3J                |            | 2.4<br>95°F            |            | 100          | <b>Sand (SP)</b><br>Similar to above, dry to damp, no odor, refusal at 2' bgl |             |
| 2           |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |
|             |              |      |                               |            |                        |            |              |   |             |



**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.968' W 107°58.694'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5443.684  
**Site Coordinates:**  
**N** 74641.537      **E** 181514.208

**Boring No.:** AOC 19-8  
**Start Date:** 8/15/2011 15:30  
**Finish Date:** 8/15/2011 15:50

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1540 | G/2V/<br>2E/3J                |            | 4.8<br>93°F            |            | 100          | Ground Surface<br><b>Silt (ML)</b><br>Very fine grain, loose, dry, light brown, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.969' W 107°58.693'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5443.599  
**Site Coordinates:**  
**N** 74643.172      **E** 181515.364

**Boring No.:** AOC-19-9  
**Start Date:** 8/16/2011 13:15  
**Finish Date:** 8/16/2011 13:30

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |  |             |
| 0           | 0-0.5'       | 1320 | G/2V/<br>2E/3J                |            | 3.2<br>91°F            |            | Ground Surface   | 0           |
|             |              |      |                               |            |                        |            | <b>Sand (SP)</b><br>Very fine grain, loose to compact, dry to damp, light gray, no odor, refusal at 0.5' bgl |             |
|             |              |      |                               |            |                        |            | Total Depth = 0.5' BGL   |             |
| 2           |              |      |                               |            |                        |            |  | 2           |
| 4           |              |      |                               |            |                        |            |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.968' W 107°58.694'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5443.81  
**Site Coordinates:**  
**N** 74644.548      **E** 181517.072

**Boring No.:** AOC 19-10  
**Start Date:** 8/15/2011 15:50  
**Finish Date:** 8/15/2011 16:10

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 1555 | G/2V/<br>2E/3J                |            | 2.6<br>93°F            |            | 100          | Ground Surface<br><b>Sand (SP)</b><br>Fine grain, loose, dry, very light gray, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates: N 36°41.975' W 107°58.704'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5439.073  
**Site Coordinates:**  
**N** 74680.844      **E** 181466.492

**Boring No.:** AOC 19-11  
**Start Date:** 8/15/2011 16:10  
**Finish Date:** 8/15/2011 16:20

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1615 | G/2V/<br>2E/3J                |            | 3.1<br>93°F            |            | 100          | Ground Surface<br><b>Silt (ML)</b><br>Very fine grain, compact, dry, light brown, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.976' W 107°58.704'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5439.119  
**Site Coordinates:**  
**N** 74687.795      **E** 181474.856

**Boring No.:** AOC 19-12  
**Start Date:** 8/15/2011 16:25  
**Finish Date:** 8/15/2011 16:35

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 1630 | G/2V/<br>2E/3J                |            | 2.8<br>92°F            |            | 100          | Ground Surface<br><b>Silty Sand (SM)</b><br>Fine grain, loose, dry, brown, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.979' W 107°58.704'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5437.428  
**Site Coordinates:**  
**N** 74702.828      **E** 181462.667

**Boring No.:** AOC 19-13  
**Start Date:** 8/15/2011 16:40  
**Finish Date:** 8/15/2011 16:50

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |  |             |
| 0           | 0-0.5'       | 1645 | G/2V/<br>2E/3J                |            | 1.8<br>92°F            |            | Ground Surface   | 0           |
|             |              |      |                               |            |                        |            | <b>Silty Sand (SM)</b><br>Very fine grain, loose, dry, light gray, no odor, detritus |             |
|             |              |      |                               |            |                        |            | Total Depth = 0.5' BGL   |             |
| 2           |              |      |                               |            |                        |            |  | 2           |
| 4           |              |      |                               |            |                        |            |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.969' W 107°58.700'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5448.15  
**Site Coordinates:**  
**N** 74631.436 **E** 181457.65

**Boring No.:** AOC 19-14  
**Start Date:** 8/16/2011 08:00  
**Finish Date:** 8/16/2011 08:20

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 0810 | G/2V/<br>2E/3J                |            | 2.4<br>66°F            |            | 100          | Ground Surface<br><br><b>Silt (ML)</b><br>Very fine grain, loose, dry, brown, no odor | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.968' W 107°58.701'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5445.7  
**Site Coordinates:**  
**N** 74629.846      **E** 181451.326

**Boring No.:** AOC 19-15  
**Start Date:** 8/16/2011 08:25  
**Finish Date:** 8/16/2011 08:40

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 0830 | G/2V/<br>2E/3J                |            | 1.2<br>67°F            |            | 100          | Ground Surface<br><b>Silt (ML)</b><br>Very fine grain, compact, dense, dry, no odor, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |



**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.967' W 107°58.702'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5447.474  
**Site Coordinates:**  
**N** 74628.77      **E** 181446.148

**Boring No.:** AOC 19-16  
**Start Date:** 8/16/2011 08:40  
**Finish Date:** 8/16/2011 09:00

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 0850 | G/2V/<br>2E/3J                |            | 4.6<br>69°F            |            | 100          | Ground Surface<br><b>Silty Sand (SM)</b><br>Fine grain, loose, dry, tan, no odor, gravel present, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.966' W 107°58.699'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5446.746  
**Site Coordinates:**  
**N** 74622.64      **E** 181451.735

**Boring No.:** AOC 19-17  
**Start Date:** 8/16/2011 09:00  
**Finish Date:** 8/16/2011 09:20

| Depth (ft.) | Sampling        |      |                               |            |                        |            | Sample Description   | Depth (ft.) |
|-------------|-----------------|------|-------------------------------|------------|------------------------|------------|--|-------------|
|             | Sample Depth    | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |  |             |
| 0           | 0-0.5'<br>& Dup | 0910 | G/4V/<br>4E/6J                |            | 3.2<br>69°F            |            | Ground Surface   | 0           |
|             |                 |      |                               |            |                        |            | <b>Sandy Clay (CL)</b><br>Low plasticity, soft, damp to moist, dark brown, no odor, roots, loamy with earthworms, detritus |             |
|             |                 |      |                               |            |                        |            | Total Depth = 0.5' BGL   |             |
| 2           |                 |      |                               |            |                        |            |  | 2           |
| 4           |                 |      |                               |            |                        |            |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.958' W 107°58.696'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5460.765  
**Site Coordinates:**  
**N** 74575.084      **E** 181497.765

**Boring No.:** AOC 19-18  
**Start Date:** 8/16/2011 13:30  
**Finish Date:** 8/16/2011 13:50

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1345 | G/2V/<br>2E/3J                |            | 4.2<br>91°F            |            | 100          | Ground Surface<br><b>Gravelly Sand (SW)</b><br>Fine to medium grain, loose to compact, dry, brown, no odor, roots/stems/bark present, refusal at 0.5' bgl in large gravel | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.958' W 107°58.695'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5461.075  
**Site Coordinates:**  
**N** 74575.963      **E** 181503.283

**Boring No.:** AOC 19-19  
**Start Date:** 8/16/2011 11:00  
**Finish Date:** 8/16/2011 11:15

| Depth (ft.) | Sampling     |      |                               |            |                        | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) |              |  |             |
| 0           | 0-0.5'       | 1105 | G/2V/<br>2E/3J                |            | 2.8<br>83°F            | 100          | Ground Surface<br><br><b>Gravelly Sand (SW)</b><br>Very fine grain, loose, dry, tan, no odor, gravel ranged from 1/2 inch to +3 inches, subrounded, detritus | 0           |
| 2           |              |      |                               |            |                        |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |              |  | 4           |

**RPS****LOG OF BORING****Client:** Western Refining Southwest, Inc.**Site:** SWMU Group #6, Bloomfield Refinery**Job No.:** UEC01318 - Bloomfield, NM**Geologist:** Tracy Payne**Driller:** N/A**Drilling Rig:** N/A**Drilling Method:** N/A**Sampling Method:** Stainless Steel Trowel**Comments:** GPS Coordinates - N 36°41.950' W 107°58.698'**Total Depth:** 0.5' bgl**Ground Water:** Not Encountered**Elev., TOC (ft. msl):** --**Elev., PAD (ft. msl):** --**Elev., GL (ft. msl):** 5477.244**Site Coordinates:****N** 74532.91**E** 181482.057**Boring No.:** AOC 19-20**Start Date:** 8/16/2011 10:30**Finish Date:** 8/16/2011 10:55

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 1045 | G/2V/<br>2E/3J                |            | 3.2<br>82°F            |            | 100          | Ground Surface<br><b>Gravelly Sand (SW)</b><br>Very fine grain, loose, dry, tan, no odor, gravel ranged from 1/2 inch to +3 inches, subrounded, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.958' W 107°58.697'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5461.77  
**Site Coordinates:**  
N 74574.618 E 181492.695

**Boring No.:** AOC 19-21  
**Start Date:** 8/16/2011 10:00  
**Finish Date:** 8/16/2011 10:10

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 1005 | G/2V/<br>2E/3J                |            | 3.6<br>80°F            |            | 100          | Ground Surface<br><b>Gravelly Sand (SW)</b><br>Very fine grain, loose, dry, tan, no odor, gravel ranged from 1/2 inch to +3 inches, subrounded, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.949' W 107°58.696'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5477.52  
**Site Coordinates:**  
**N** 74530.609      **E** 181493.6

**Boring No.:** AOC 19-22  
**Start Date:** 8/16/2011 10:20  
**Finish Date:** 8/16/2011 10:30

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 1025 | G/2V/<br>2E/3J                |            | 2.1<br>80°F            |            | 100          | Ground Surface<br><br><b>Gravelly Sand (SW)</b><br>Very fine grain, loose, dry, tan, no odor, gravel ranged from 1/2 inch to +3 inches, subrounded, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.947' W 107°58.699'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5480.975  
**Site Coordinates:**  
**N** 74513.176      **E** 181482.442

**Boring No.:** AOC 19-23  
**Start Date:** 8/16/2011 09:30  
**Finish Date:** 8/16/2011 09:55

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 0945 | G/2V/<br>2E/3J                |            | 4.9<br>77°F            |            | 100          | Ground Surface<br><b>Weathered Sandstone (SS)</b><br>Fine grain, moderate to poor cementation, dry, dark brown, stained, faint odor | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |




**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.939' W 107°58.740'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5469.417  
**Site Coordinates:**  
**N** 74461.182      **E** 181283.544

**Boring No.:** AOC 20-1**Start Date:** 8/17/2011 07:45**Finish Date:** 8/17/2011 08:10

| Depth (ft.) | Sampling     |      |                               |            |                        |  | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|--|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class   |              |  |             |
| 0           | 0-0.5'       | 0800 | G/2V/<br>2E/3J                |            | 2.0<br>67°F            |  | 100          | Ground Surface<br><br><b>Gravelly Sand (SW)</b><br>Fine grain, loose, damp, brown, no odor, gravel ranges from 1/2 inch to +2 inches, subrounded | 0           |
| 2           |              |      |                               |            |                        |  |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |  |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.940' W 107°58.738'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5467.509  
**Site Coordinates:**  
**N** 74472.531      **E** 181289.678

**Boring No.:** AOC 20-2**Start Date:** 8/17/2011 08:15**Finish Date:** 8/17/2011 08:30

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 0820 | G/2V/<br>2E/3J                |            | 1.3<br>70°F            |            | 100          | Ground Surface<br><b>Clayey Sand (SC)</b><br>Fine grain, compact, damp, brown, no odor | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.940' W 107°58.739'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5465.067  
**Site Coordinates:**  
**N** 74471.467      **E** 181280.71

**Boring No.:** AOC 20-3  
**Start Date:** 8/17/2011 08:35  
**Finish Date:** 8/17/2011 08:45

| Depth (ft.) | Sampling        |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|-----------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth    | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'<br>& Dup | 0840 | G/4V/<br>4E/6J                |            | 33<br>71°F             |            | 100          | Ground Surface<br><br>Clay (CL)<br>Low plasticity, soft, damp, tan and black, faint odor | 0           |
| 2           |                 |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   | 2           |
| 4           |                 |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.939' W 107°58.742'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5467.38  
**Site Coordinates:**  
**N** 74470.728      **E** 181275.869

**Boring No.:** AOC 20-4  
**Start Date:** 8/17/2011 08:45  
**Finish Date:** 8/17/2011 09:05

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 0900 | G/2V/<br>2E/3J                |            | 9.4<br>74°F            |            | 100          | Ground Surface   | 0           |
|             |              |      |                               |            |                        |            |              | <b>Silt (ML)</b><br>Low plasticity, soft, damp, brown, no odor |             |
|             |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   |             |
| 2           |              |      |                               |            |                        |            |              |  | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.948' W 107°58.739'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5456.308  
**Site Coordinates:**  
**N** 74513.196      **E** 181276.971

**Boring No.:** AOC 20-5  
**Start Date:** 8/17/2011 09:15  
**Finish Date:** 8/17/2011 10:00

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |  |             |
| 0           | 0-0.5'       | 0945 | G/2V/<br>2E/3J                |            | 2.4<br>80°F            |            | Ground Surface   | 0           |
|             |              |      |                               |            |                        |            | <b>Sandy Gravel (GP)</b><br>Rounded gravel, 1/2 inch to +6 inches, very fine grain sand, loose, dry, no odor, brown, refusal at 0.5' bgl |             |
|             |              |      |                               |            |                        |            | Total Depth = 0.5' BGL   |             |
| 2           |              |      |                               |            |                        |            |  | 2           |
| 4           |              |      |                               |            |                        |            |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.950' W 107°58.740'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5453.998  
**Site Coordinates:**  
**N** 74526.439      **E** 181278.604

**Boring No.:** AOC 20-6  
**Start Date:** 8/17/2011 10:00  
**Finish Date:** 8/17/2011 10:10

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1005 | G/2V/<br>2E/3J                |            | 2.4<br>83°F            |            | 100          | Ground Surface<br><br><b>Sand (SP)</b><br>Very fine grain, loose, dry, brown, no odor | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.950' W 107°58.743'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5453.359  
**Site Coordinates:**  
**N** 74530.099      **E** 181266.994

**Boring No.:** AOC 20-7  
**Start Date:** 8/17/2011 10:15  
**Finish Date:** 8/17/2011 10:30

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1025 | G/2V/<br>2E/3J                |            | 0.7<br>84°F            |            | 100          | Ground Surface<br><br><b>Sand (SP)</b><br>Fine grain, loose, dry, light gray, no odor | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.951' W 107°58.742'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5451.499  
**Site Coordinates:**  
**N** 74535.077      **E** 181273.44

**Boring No.:** AOC 20-8  
**Start Date:** 8/17/2011 10:35  
**Finish Date:** 8/17/2011 10:50

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |  |             |
| 0           | 0-0.5'       | 1045 | G/2V/<br>2E/3J                |            | 0.7<br>84°F            |            | Ground Surface   | 0           |
|             |              |      |                               |            |                        |            | <b>Sandy Gravel (GP)</b><br>Subrounded, 1/2 inch to +6 inches, fine grain sand, compact, dry, brown, no odor |             |
|             |              |      |                               |            |                        |            | Total Depth = 0.5' BGL   |             |
| 2           |              |      |                               |            |                        |            |  | 2           |
| 4           |              |      |                               |            |                        |            |  | 4           |



**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.951' W 107°58.739'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5451.334  
**Site Coordinates:**  
**N** 74537.171      **E** 181282.135

**Boring No.:** AOC 20-9  
**Start Date:** 8/17/2011 11:00  
**Finish Date:** 8/17/2011 11:30

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 1120 | G/2V/<br>2E/3J                |            | 2.4<br>95°F            |            | 100          | Ground Surface   | 0           |
|             |              |      |                               |            |                        |            |              | <b>Sand (SP)</b><br>Very fine grain, loose, dry, light gray, no odor |             |
|             |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL   |             |
| 2           |              |      |                               |            |                        |            |              |  | 2           |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.951' W 107°58.743'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5450.617  
**Site Coordinates:**  
**N** 74540.095      **E** 181268.539

**Boring No.:** AOC 20-10  
**Start Date:** 8/17/2011 11:30  
**Finish Date:** 8/17/2011 11:50

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1140 | G/2V/<br>2E/3J                |            | 2.7<br>92°F            |            | 100          | Ground Surface<br><br><b>Sand (SP)</b><br>Very fine grain, loose, dry, light tan, no odor | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING****Client:** Western Refining Southwest, Inc.**Site:** SWMU Group #6, Bloomfield Refinery**Job No.:** UEC01318 - Bloomfield, NM**Geologist:** Tracy Payne**Driller:** N/A**Drilling Rig:** N/A**Drilling Method:** N/A**Sampling Method:** Stainless Steel Trowel**Comments:** GPS Coordinates - N 36°41.951' W 107°58.742'**Total Depth:** 0.5' bgl**Ground Water:** Not Encountered**Elev., TOC (ft. msl):** --**Elev., PAD (ft. msl):** --**Elev., GL (ft. msl):** 5449.867**Site Coordinates:****N** 74540.076**E** 181274.717**Boring No.:** AOC 20-11**Start Date:** 8/17/2011 11:55**Finish Date:** 8/17/2011 12:10

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1200 | G/2V/<br>2E/3J                |            | 2.3<br>94°F            |            | 100          | Ground Surface<br><b>Sandy Gravel (GP)</b><br>Subrounded, 1/2 inch to +6 inches, very fine grain sand, compact, dry, brown, no odor | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.954' W 107°58.742'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5446.088  
**Site Coordinates:**  
**N** 74555.04      **E** 181272.202

**Boring No.:** AOC 20-12  
**Start Date:** 8/17/2011 12:10  
**Finish Date:** 8/17/2011 12:25

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |   |             |
| 0           | 0-0.5'       | 1220 | G/2V/<br>2E/3J                |            | 2.1<br>93°F            |            | Ground Surface  | 0           |
|             |              |      |                               |            |                        |            | <b>Sand (SP)</b><br>Very fine grain, loose, dry, light tan, no odor |             |
|             |              |      |                               |            |                        |            | Total Depth = 0.5' BGL  |             |
| 2           |              |      |                               |            |                        |            |   | 2           |
| 4           |              |      |                               |            |                        |            |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.937' W 107°58.788'

**Total Depth:** 2.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5473.066  
**Site Coordinates:**  
**N** 74450.302      **E** 181046.206

**Boring No.:** AOC 21-1**Start Date:** 8/17/2011 14:30**Finish Date:** 8/17/2011 15:00

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |  |             |
| 0           | 0-0.5'       | 1435 | G/2V/<br>2E/3J                |            | 1.5<br>92°F            |            | 100          | Ground Surface   | 0           |
|             |              |      |                               |            |                        |            |              | <b>Silty Sand (SM)</b><br>Very fine grain, loose, dry, gray, no odor     |             |
|             | 1.5-2'       | 1445 | G/2V/<br>2E/3J                |            | 0.8<br>92°F            |            | 100          | <b>Silty Sand (SM)</b><br>Similar to above, no odor                      |             |
|             |              |      |                               |            |                        |            |              | <b>Silty Sand (SM)</b><br>Similar to above, no odor                      |             |
| 2           | 2-2.5'       | 1455 | G/2V/<br>2E/3J                |            | 1.6<br>92°F            |            | 100          | <b>Silty Sand (SM)</b><br>Similar to above, no odor, refusal at 2.5' bgl | 2           |
|             |              |      |                               |            |                        |            |              | <b>Silty Sand (SM)</b><br>Similar to above, no odor, refusal at 2.5' bgl |             |
|             |              |      |                               |            |                        |            |              | Total Depth = 2.5' BGL   |             |
| 4           |              |      |                               |            |                        |            |              |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.940' W 107°58.786'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5472.78  
**Site Coordinates:**  
N 74460.204 E 181050.86

**Boring No.:** AOC 21-2  
**Start Date:** 8/15/2011 09:30  
**Finish Date:** 8/15/2011 09:55

| Depth (ft.) | Sampling        |      |                               |            |                        |            | Sample Description   | Depth (ft.) |
|-------------|-----------------|------|-------------------------------|------------|------------------------|------------|--|-------------|
|             | Sample Depth    | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |  |             |
| 0           | 0-0.5'<br>& Dup | 0940 | G/4V/<br>4E/6J                |            | 9.8<br>76°F            |            | Ground Surface   | 0           |
|             |                 |      |                               |            |                        |            | <b>Silty Clay (CL)</b><br>Low plasticity, soft, dry, light tan to dark gray, no odor, detritus |             |
|             |                 |      |                               |            |                        |            | Total Depth = 0.5' BGL   |             |
| 2           |                 |      |                               |            |                        |            |  | 2           |
| 4           |                 |      |                               |            |                        |            |  | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.938' W 107°58.789'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5471.977  
**Site Coordinates:**  
**N** 74465.222      **E** 181040.544

**Boring No.:** AOC 21-3  
**Start Date:** 8/15/2011 09:56  
**Finish Date:** 8/15/2011 10:11

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1000 | G/2V/<br>2E/3J                |            | 1.3<br>78°F            |            | 100          | Ground Surface<br><br><b>Silty Sand (SM)</b><br>Very fine grain, loose, dry, light tan to light gray, no odor, sandstone gravel present, detritus | 0           |
| 2           |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |

**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.939' W 107°58.788'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5471.458  
**Site Coordinates:**  
**N** 74460.782      **E** 181045.136

**Boring No.:** AOC 21-4  
**Start Date:** 8/15/2011 10:12  
**Finish Date:** 8/15/2011 10:25

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Recovery (%) | Sample Description  | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--------------|---|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |              |   |             |
| 0           | 0-0.5'       | 1020 | G/2V/<br>2E/3J                |            | 5.0<br>79°F            |            | 100          | Ground Surface  | 0           |
|             |              |      |                               |            |                        |            |              | <b>Silt (ML)</b><br>Very fine grain, loose, dry, light gray, no odor, roots and stems present, detritus |             |
|             |              |      |                               |            |                        |            |              | Total Depth = 0.5' BGL  |             |
| 2           |              |      |                               |            |                        |            |              |   | 2           |
| 4           |              |      |                               |            |                        |            |              |   | 4           |



**RPS****LOG OF BORING**

**Client:** Western Refining Southwest, Inc.  
**Site:** SWMU Group #6, Bloomfield Refinery  
**Job No.:** UEC01318 - Bloomfield, NM  
**Geologist:** Tracy Payne  
**Driller:** N/A  
**Drilling Rig:** N/A  
**Drilling Method:** N/A  
**Sampling Method:** Stainless Steel Trowel  
**Comments:** GPS Coordinates - N 36°41.940' W 107°58.789'

**Total Depth:** 0.5' bgl  
**Ground Water:** Not Encountered  
**Elev., TOC (ft. msl):** --  
**Elev., PAD (ft. msl):** --  
**Elev., GL (ft. msl):** 5470.482  
**Site Coordinates:**  
**N** 74472.412      **E** 181047.834

**Boring No.:** AOC 21-5  
**Start Date:** 8/17/2011 15:00  
**Finish Date:** 8/17/2011 16:05

| Depth (ft.) | Sampling     |      |                               |            |                        |            | Sample Description   | Depth (ft.) |
|-------------|--------------|------|-------------------------------|------------|------------------------|------------|--|-------------|
|             | Sample Depth | Time | Sample Type/<br>Container/No. | Saturation | Organic Vapor<br>(ppm) | USCS Class |  |             |
| 0           | 0-0.5'       | 1520 | G/2V/<br>2E/3J                |            | 0.7<br>91°F            |            | Ground Surface   | 0           |
|             |              |      |                               |            |                        |            | <b>Silty Sand (SM)</b><br>Very fine grain, loose, dry, gray, no odor |             |
|             | 1.5-2'       | 1540 | G/2V/<br>2E/3J                |            | 0.7<br>91°F            |            | <b>Silty Sand (SM)</b><br>Similar to above, no odor                  |             |
|             |              |      |                               |            |                        |            |  |             |
| 2           | 2-2.5'       | 1600 | G/2V/<br>2E/3J                |            | 0.7<br>91°F            |            | <b>Silty Sand (SM)</b><br>Similar to above, no odor                  | 2           |
|             |              |      |                               |            |                        |            | Similar to above, no odor, refusal at 2.5' bgl                       |             |
|             |              |      |                               |            |                        |            | Total Depth = 2.5' BGL   |             |
| 4           |              |      |                               |            |                        |            |  | 4           |

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

The DAF value was calculated using equation 19 from NMED's *Technical Background Document for Development of Soil Screening Levels (Revision 5.0, August 2009)*.

$$DAF = 1 + \left( \frac{K * i * D}{I * L} \right) \quad DAF = 1 + \left( \frac{4,893 * 0.0023 * 1}{0.01 * 100} \right) = 11.25$$

Where:

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

- 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 = 4,893 m/yr as determined from pumping test at well RW-22 (lowest of three values determined during 1994 RCRA Facility Investigation)
- i = 0.0023 m/m as measured during August 2008 ground water sampling event
- D = 1 m (lower of aquifer thickness (1m) or calculated mixing zone depth (10.58m))
- I = 0.01 m derivation using EPA's HELP model as described below
- L = 100 m – conservative average of SWMU/AOC source area length
- D<sub>a</sub> = 1 m - average saturated thickness measured during August 2008 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 Bloomfield 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 (Bloomfield 3 SE) as part of the NWS Cooperator Climate Stations. The weather station is located 1.7 miles south of Bloomfield on HWY 44 and then two miles east on Industrial Blvd, thus being approximately two miles southeast of the Western Bloomfield Refinery.

Data obtained from the Bloomfield 3 SE station includes mean monthly temperature and average monthly precipitation. The average wind speed (13.5 km/hr) was obtained from

the Western Regional Climate Center, as measured at the Farmington, 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 Bloomfield 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 Bloomfield. The quarterly relative humidity values used are 48%, 30%, 45%, and 50% for the first, second, third, and fourth quarters, respectively.

The vadose zone physical properties were based on the predominant lithology as observed during on-site monitoring well installation. The soil type chosen in the model was loamy sand with an average thickness of 5 meters. The land surface was assumed to be bare soil with a slope of 0%. This should be a conservative estimate, as there is a slight slope across most of the refinery with the exception of areas within tank dikes. There are structures (e.g., parking lots, building pads, concrete foundations, etc.) that could limit infiltration but the model assumes only bare soil without any obstructions to infiltration. Based on the selected soil type (loamy sand), the model default value for porosity is 0.437, field capacity is 0.105, wilting point is 0.047, and saturated hydraulic conductivity is 0.0017 cm/day. These model default values are taken from the US Department of Agriculture.

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). The model output is enclosed, showing the annual values. Over the modeled 40 year period, the average annual infiltration was 0.01 meters. This average annual infiltration was used in the aforementioned calculation of the site-specific DAF value.

### Bloomfield 3 SE, New Mexico Weather Station Data

|                          | Jan.  | Feb.   | March  | April | May    | June  | July   | August | Sept. | Oct.  | Nov.   | Dec.   | Annual |
|--------------------------|-------|--------|--------|-------|--------|-------|--------|--------|-------|-------|--------|--------|--------|
| Average Max Temp (F)     | 41    | 48.6   | 57.4   | 67.2  | 77.4   | 88    | 92     | 89     | 81.8  | 69.4  | 54.4   | 43.4   | 67.5   |
| Average Min. Temp (F)    | 16.2  | 22.4   | 27.8   | 35    | 43.8   | 52.1  | 59.6   | 57.7   | 49.5  | 37.7  | 25.7   | 18     | 37.1   |
| Mean Monthly Temp (F)    | 28.6  | 35.5   | 42.6   | 51.1  | 60.6   | 70.1  | 75.8   | 73.4   | 65.7  | 53.6  | 40.1   | 30.7   | 52.3   |
| Mean Monthly Temp (C)    | -1.89 | 1.94   | 5.89   | 10.61 | 15.89  | 21.14 | 24.33  | 22.97  | 18.69 | 11.97 | 4.47   | -0.72  |        |
| Average total Prec. (in) | 0.55  | 0.56   | 0.63   | 0.6   | 0.52   | 0.38  | 0.99   | 1.27   | 0.95  | 0.95  | 0.63   | 0.57   | 8.60   |
| Average total Prec. (mm) | 13.97 | 14.224 | 16.002 | 15.24 | 13.208 | 9.652 | 25.146 | 32.258 | 24.13 | 24.13 | 16.002 | 14.478 |        |

Data collected from 1/1/1914 to 12/31/2005 at the Bloomfield 3 SE (#291063) weather monitoring station; obtained from Western Regional Climate Center, National Oceanic & Atmospheric Administration

### Albuquerque, New Mexico Weather Station Data

|                          | Jan.  | Feb.  | March  | April  | May   | June   | July  | August | Sept.  | Oct.   | Nov.   | Dec.   | Annual |
|--------------------------|-------|-------|--------|--------|-------|--------|-------|--------|--------|--------|--------|--------|--------|
| Average Max Temp (F)     | 49.9  | 53.6  | 60.8   | 72.4   | 80.1  | 91.1   | 93.2  | 92     | 84.8   | 74.4   | 59.5   | 49.8   | 71.8   |
| Average Min. Temp (F)    | 20.9  | 23.7  | 30.2   | 38.7   | 46.8  | 56.3   | 62    | 60.5   | 52.5   | 40.3   | 26.4   | 21.6   | 40     |
| Mean Monthly Temp (F)    | 35.4  | 38.65 | 45.5   | 55.55  | 63.45 | 73.7   | 77.6  | 76.25  | 68.65  | 57.35  | 42.95  | 35.7   | 55.9   |
| Mean Monthly Temp (C)    | 1.89  | 3.69  | 7.50   | 13.08  | 17.47 | 23.17  | 25.33 | 24.58  | 20.36  | 14.08  | 6.08   | 2.06   |        |
| Average total Prec. (in) | 0.32  | 0.29  | 0.46   | 0.61   | 0.7   | 0.87   | 1.3   | 1.57   | 1.03   | 0.63   | 0.43   | 0.46   | 8.67   |
| Average total Prec. (mm) | 8.128 | 7.366 | 11.684 | 15.494 | 17.78 | 22.098 | 33.02 | 39.878 | 26.162 | 16.002 | 10.922 | 11.684 |        |

Data collected from 1/1/1932 to 1/31/1954 at the Albuquerque (#290222) weather monitoring station; obtained from Western Regional Climate Center, National Oceanic & Atmospheric Administration

F - Fahrenheit  
C - Celsius  
in - inch  
mm - millimeter

# ***Project : Western Refining Bloomfield, New Mexico***

*Estimation of infiltration at Bloomfield Refinery*

***Model : HELP***

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

***Author : Scott Crouch***

***Client : Western Refining - Randy Schmaltz***

***Location : Bloomfield, NM***

***3/11/2010***

## Profile 1

### 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 | Bare soil | (-)   |

### Profile Structure

| Layer  | Top ( m) | Bottom ( m) | Thickness ( m) |
|--|----------|-------------|----------------|
|  Loamy Sand | 100.0000 | 95.0000     | 5.0000         |

#### 1.1. Layer. Loamy Sand

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

[HELP] Vertical Perc. Layer Parameters

| Parameter             | Value  | Units     |
|-----------------------|--------|-----------|
| total porosity        | 0.437  | (vol/vol) |
| field capacity        | 0.105  | (vol/vol) |
| wilting point         | 0.047  | (vol/vol) |
| sat.hydr.conductivity | 0.0017 | (cm/sec)  |
| subsurface inflow     | 0      | (mm/year) |

# Annual Totals rate (m)

|             | Precipitation (m) | Runoff (m) | Evapotranspiration (m) | Percolation or leakage through Layer 1 (m) |
|-------------|-------------------|------------|------------------------|--|
| Year-1 (m)  | 1.9660E-01        | 0.0000E+00 | 1.8579E-01             | 5.2109E-05                                 |
| Year-2 (m)  | 3.0180E-01        | 0.0000E+00 | 2.6922E-01             | 1.0255E-04                                 |
| Year-3 (m)  | 2.3510E-01        | 0.0000E+00 | 2.3452E-01             | 1.9650E-04                                 |
| Year-4 (m)  | 2.3000E-01        | 0.0000E+00 | 2.1004E-01             | 2.4626E-04                                 |
| Year-5 (m)  | 2.5270E-01        | 0.0000E+00 | 2.3977E-01             | 4.1142E-04                                 |
| Year-6 (m)  | 1.5870E-01        | 0.0000E+00 | 1.4899E-01             | 3.6109E-04                                 |
| Year-7 (m)  | 1.8420E-01        | 0.0000E+00 | 1.7010E-01             | 5.0670E-04                                 |
| Year-8 (m)  | 2.5770E-01        | 0.0000E+00 | 2.3978E-01             | 5.9778E-04                                 |
| Year-9 (m)  | 1.9170E-01        | 0.0000E+00 | 1.7956E-01             | 7.2288E-04                                 |
| Year-10 (m) | 2.2820E-01        | 0.0000E+00 | 1.9825E-01             | 9.4104E-04                                 |
| Year-11 (m) | 2.3680E-01        | 0.0000E+00 | 2.2456E-01             | 1.6311E-03                                 |
| Year-12 (m) | 2.5940E-01        | 0.0000E+00 | 2.4152E-01             | 3.7601E-03                                 |
| Year-13 (m) | 1.8440E-01        | 0.0000E+00 | 1.7107E-01             | 5.6153E-03                                 |
| Year-14 (m) | 1.5860E-01        | 0.0000E+00 | 1.5145E-01             | 1.0341E-02                                 |
| Year-15 (m) | 2.4990E-01        | 0.0000E+00 | 2.3436E-01             | 1.4166E-02                                 |
| Year-16 (m) | 1.6700E-01        | 0.0000E+00 | 1.5633E-01             | 1.4482E-02                                 |
| Year-17 (m) | 1.3040E-01        | 0.0000E+00 | 1.1372E-01             | 1.2954E-02                                 |
| Year-18 (m) | 1.5020E-01        | 0.0000E+00 | 1.4066E-01             | 1.3977E-02                                 |
| Year-19 (m) | 2.0530E-01        | 0.0000E+00 | 1.9662E-01             | 1.3219E-02                                 |
| Year-20 (m) | 1.8180E-01        | 0.0000E+00 | 1.6946E-01             | 1.0024E-02                                 |
| Year-21 (m) | 2.3550E-01        | 0.0000E+00 | 2.1477E-01             | 1.0887E-02                                 |
| Year-22 (m) | 1.3750E-01        | 0.0000E+00 | 1.3022E-01             | 1.0618E-02                                 |
| Year-23 (m) | 2.3340E-01        | 0.0000E+00 | 2.2529E-01             | 1.4634E-02                                 |
| Year-24 (m) | 2.2170E-01        | 0.0000E+00 | 2.0414E-01             | 1.0021E-02                                 |
| Year-25 (m) | 1.4510E-01        | 0.0000E+00 | 1.3452E-01             | 1.3558E-02                                 |
| Year-26 (m) | 2.0130E-01        | 1.2902E-06 | 1.7333E-01             | 1.3059E-02                                 |
| Year-27 (m) | 2.3200E-01        | 0.0000E+00 | 2.1409E-01             | 1.5689E-02                                 |
| Year-28 (m) | 1.9260E-01        | 0.0000E+00 | 1.8730E-01             | 9.9471E-03                                 |
| Year-29 (m) | 2.3390E-01        | 0.0000E+00 | 2.1475E-01             | 1.1847E-02                                 |
| Year-30 (m) | 1.8890E-01        | 0.0000E+00 | 1.7801E-01             | 1.8487E-02                                 |
| Year-31 (m) | 2.4520E-01        | 0.0000E+00 | 2.2175E-01             | 1.6094E-02                                 |
| Year-32 (m) | 2.2790E-01        | 0.0000E+00 | 2.0877E-01             | 1.2385E-02                                 |
| Year-33 (m) | 3.1730E-01        | 4.0020E-04 | 2.9335E-01             | 1.3069E-02                                 |
| Year-34 (m) | 2.1170E-01        | 0.0000E+00 | 1.8598E-01             | 1.4984E-02                                 |
| Year-35 (m) | 2.7430E-01        | 0.0000E+00 | 2.6796E-01             | 1.6877E-02                                 |
| Year-36 (m) | 1.5090E-01        | 0.0000E+00 | 1.2899E-01             | 2.4361E-02                                 |
| Year-37 (m) | 2.1680E-01        | 0.0000E+00 | 2.1801E-01             | 1.6731E-02                                 |
| Year-38 (m) | 1.7490E-01        | 0.0000E+00 | 1.5227E-01             | 1.8959E-02                                 |
| Year-39 (m) | 2.1190E-01        | 0.0000E+00 | 1.6801E-01             | 1.5479E-02                                 |
| Year-40 (m) | 1.7540E-01        | 0.0000E+00 | 1.8233E-01             | 1.7584E-02                                 |
| Total (m)   | 8.3887E+00        | 4.0149E-04 | 7.7796E+00             | 3.9958E-01                                 |

Average  
= 0.01/m



# NWS Cooperator CLIMATE STATIONS

WEATHER DATA FROM INDIVIDUAL STATIONS AROUND  
NEW MEXICO

## Bloomfield 3-SE-Bloomfield, NM

### Climate Data

NO PICTURE  
AVAILABLE

Location:  
From  
Bloomfield,  
NM go 1.7  
miles south  
on HWY 44,  
turn east on  
Industrial  
BLVD and  
go 2.0 miles  
to gas  
compressor  
plant on  
right.

Elevation:  
5806 feet

Latitude:  
36°40'

Longitude:  
107°58'

Ground Cover: Flat sandy plateau cut by broken terrain of sandstone hills and arroyos.

Cooperator Number: 29-1063-1

Questions or comments about this page can be directed to:



[webmaster@weather.nmsu.edu](mailto:webmaster@weather.nmsu.edu)  
NMSU Weather BBS  
Dept. of Agronomy and Horticulture  
BOX 30003, Dept. 3Q  
LAS CRUCES, NM 88003-0003

NMSU MONITORED CLIMATE  
STATIONS  
NMSU Weather Homepage

# Western Regional Climate Center

## About Us

The Regional Climate Centers (RCC) deliver climate services at national, regional and state levels working with NOAA partners in the National Climatic Data Center, National Weather Service, the American Association of State Climatologists, and NOAA Research Institutes. This successful effort resulted in jointly developed products, services, and capabilities that enhance the delivery of climate information to the American public, and builds a solid foundation for a National Climate Service. As NOAA and Congress work to help society adapt to climate change, these collaborative efforts form a framework for the service, data stewardship, and applied research components of the National Climate Service.

[Climate Information](#) [Current Observations](#) [Projects](#) [Monitoring](#) [More Info](#) [Educational & Travel](#) [More Sources](#) [About Us](#) [BACK TO HOME PAGE](#)

---

- [Overview of the Western Regional Climate Center](#)
- [Roles of the Regional Climate Centers in Climate Services](#)
- [Pricing and Formats](#)
- [DRI Home Page](#)
- [Weather Info \(DRI ACCESS ONLY\)](#)
- [WRCC Software Library \(restricted access\)](#)
- [DRI LDM Logs \(ldmd, netcheck and syscheck\)](#)

---

Division of Atmospheric Sciences



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 |

8.4 Miles/hr x 1.60934 = 13.5 kts/hr

NEVADA

AVERAGE WIND SPEED - MPH

| STATION                | ID   | Years     | Jan | Feb | Mar | Apr  | May  | Jun  | Jul | Aug | Sep | Oct | Nov | Dec | Ann |
|------------------------|------|-----------|-----|-----|-----|------|------|------|-----|-----|-----|-----|-----|-----|-----|
| DESERT ROCK AP-MERCURY | KDRA | 1996-2006 | 8.0 | 8.8 | 9.2 | 10.7 | 10.5 | 10.5 | 9.6 | 9.1 | 8.8 | 8.2 | 7.7 | 8.4 | 9.1 |
| ELKO AIRPORT ASOS      | KEKO | 1996-2006 | 4.6 | 5.3 | 5.9 | 6.7  | 6.4  | 6.3  | 5.7 | 5.3 | 5.0 | 4.6 | 4.6 | 4.8 | 5.4 |
| ELY AIRPORT ASOS       | KELY | 1996-2006 | 9.0 | 9.0 | 9.6 | 10.3 | 9.8  | 10.2 | 9.8 | 9.9 | 9.6 | 9.5 | 8.8 | 9.2 | 9.5 |

# BLOOMFIELD 3 SE, NEW MEXICO (291063)

## 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)      | 41.0 | 48.6 | 57.4 | 67.2 | 77.4 | 88.0 | 92.0 | 89.0 | 81.8 | 69.4 | 54.4 | 43.4 | 67.5   |
| Average Min. Temperature (F)      | 16.2 | 22.4 | 27.8 | 35.0 | 43.8 | 52.1 | 59.6 | 57.7 | 49.5 | 37.7 | 25.7 | 18.0 | 37.1   |
| Average Total Precipitation (in.) | 0.55 | 0.56 | 0.63 | 0.60 | 0.52 | 0.38 | 0.99 | 1.27 | 0.95 | 0.95 | 0.63 | 0.57 | 8.61   |
| Average Total SnowFall (in.)      | 3.8  | 2.2  | 1.0  | 0.3  | 0.0  | 0.0  | 0.0  | 0.0  | 0.0  | 0.1  | 0.6  | 3.4  | 11.4   |
| 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.: 92.9% Min. Temp.: 93.2% Precipitation: 95.4% Snowfall: 79% Snow Depth: 70.2%

Check [Station Metadata](#) or [Metadata graphics](#) for more detail about data completeness.

*Western Regional Climate Center, [wrcc@dri.edu](mailto:wrcc@dri.edu)*

Back to:



NOTE:

To print data frame (right side), click on right frame before printing.

1971 - 2000

- [Daily Temp. & Precip.](#)
- [Daily Tabular data \(~23 KB\)](#)
- [Monthly Tabular data \(~1 KB\)](#)
- [NCDC 1971-2000 Normals \(~3 KB\)](#)

1961 - 1990

- [Daily Temp. & Precip.](#)
- [Daily Tabular data \(~23 KB\)](#)
- [Monthly Tabular data \(~1 KB\)](#)
- [NCDC 1961-1990 Normals \(~3 KB\)](#)

Period of Record

- [Station Metadata](#)
- [Station Metadata Graphics](#)

General Climate Summary Tables

- [Temperature](#)
- [Precipitation](#)
- [Heating Degree Days](#)
- [Cooling Degree Days](#)
- [Growing Degree Days](#)

ALBUQUERQUE, NEW MEXICO (290222)

Period of Record Monthly Climate Summary

Period of Record : 1/ 1/1932 to 1/31/1954

|                                   | Jan  | Feb  | Mar  | Apr  | May  | Jun  | Jul  | Aug  | Sep  | Oct  | Nov  | Dec  | Annual |
|-----------------------------------|------|------|------|------|------|------|------|------|------|------|------|------|--------|
| Average Max. Temperature (F)      | 49.9 | 53.6 | 60.8 | 72.4 | 80.1 | 91.1 | 93.2 | 92.0 | 84.8 | 74.4 | 59.5 | 49.8 | 71.8   |
| Average Min. Temperature (F)      | 20.9 | 23.7 | 30.2 | 38.7 | 46.8 | 56.3 | 62.0 | 60.5 | 52.5 | 40.3 | 26.4 | 21.6 | 40.0   |
| Average Total Precipitation (in.) | 0.32 | 0.29 | 0.46 | 0.61 | 0.70 | 0.87 | 1.30 | 1.57 | 1.03 | 0.63 | 0.43 | 0.46 | 8.67   |
| Average Total SnowFall (in.)      | 2.3  | 1.5  | 1.0  | 0.9  | 0.2  | 0.0  | 0.0  | 0.0  | 0.0  | 0.0  | 1.2  | 2.0  | 9.3    |
| 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.: 26.1% Min. Temp.: 26.1% Precipitation: 72.2% Snowfall: 26.1% Snow Depth: 26.1%

Check [Station Metadata](#) or [Metadata graphics](#) for more detail about data completeness.

Western Regional Climate Center, [wrcc@dri.edu](mailto:wrcc@dri.edu)

# Equation 1

## Combined Exposures to Noncarcinogenic Contaminants in Soil Residential Scenario

$$C' = \frac{THQ \times BW_c \times AT_n}{EF_r \times ED_c \left[ \left( \frac{1}{RfD_o} \times \frac{IRS_c}{10^6 \text{ mg/kg}} \right) + \left( \frac{1}{RfD_o} \times \frac{SA_c \times AF_c \times ABS}{10^6 \text{ mg/kg}} \right) + \left( \frac{1}{RFD_i} \times \frac{IRA_c}{VF_s + PEF} \right) \right]}$$

| Parameter        | Definition (units)  | NMED Default | TPH C11-C21 Aromatics |
|------------------|---|--------------|-----------------------|
| C                | Contaminant concentration (mg/kg)                             | --           | <b>1.83E+03</b>       |
| THQ              | Target hazard quotient  | 1            | 1                     |
| BW <sub>c</sub>  | Body weight, child (kg)                                       | 15           | 15                    |
| AT <sub>n</sub>  | Averaging time, noncarcinogens, ED x 365 (days)               | 2190         | 2190                  |
| EF <sub>i</sub>  | Exposure frequency, resident (days/yr)                        | 350          | 350                   |
| ED <sub>c</sub>  | Exposure duration, child (yr)                                 | 6            | 6                     |
| IRS <sub>c</sub> | Soil ingestion rate, child (mg/day)                           | 200          | 200                   |
| RfDo             | Oral reference dose (mg/kg-day)                               | --           | <b>0.03</b>           |
| SA <sub>c</sub>  | Dermal surface area, child (cm <sup>2</sup> /day)             | 2800         | 2800                  |
| AF <sub>c</sub>  | Soil adherence factor, child (mg/cm <sup>2</sup> )            | 0.2          | 0.2                   |
| ABS              | Skin absorption factor (unitless)                             | --           | <b>0.1</b>            |
| IRA <sub>c</sub> | Inhalation rate, child (m <sup>3</sup> /day)                  | 10           | 10                    |
| RfDi             | Inhalation reference dose (mg/kg-day)                         | --           | <b>0.05</b>           |
| VFs              | Volatilization factor, Equation 14 (mg <sup>3</sup> /kg)      | --           | <b>8.10E+04</b>       |
| PEF              | Particulate emission factor, Equation 16 (m <sup>3</sup> /kg) | --           | <b>6.61E+09</b>       |

| Ingestion   | Dermal      | Inhalation |
|-------------|-------------|------------|
| 2346.428571 | 8380.102041 | 5.17E+08   |

Equations and Default values from New Mexico Environment Department's Technical Background Document for Development of Soil Screening Levels, Revisions 5.0, August 2009  
Toxicity factors from Massachusetts Department of Environmental Protection Massachusetts Contingency Plan Standards Spreadsheets 2009 (**bolded values**)

# Equation 14

## Derivation of the Volatilization Factor for Residential and Commercial/Industrial Scenarios

$$VF_s = \frac{Q/C_{vol} \times (3.14 \times D_A \times T)^{0.5} \times 10^{-4}}{(2 \times \rho_b \times D_A)}$$

$$D_A = \frac{\left[ \frac{\Theta_a^{10/3} D_a H' + \Theta_w^{10/3} D_w}{n^2} \right]}{\rho_b K_d + \Theta_w + \Theta_a H'}$$

| Parameter          | Definition (units)   | NMED<br>Default | TPH C11-C21<br>Aromatics |
|--------------------|--|-----------------|--------------------------|
| VF <sub>s</sub>    | Volatilization factor for soil (m <sup>3</sup> /kg)  | --              | <b>8.10E+04</b>          |
| D <sub>A</sub>     | Apparent diffusivity (cm <sup>2</sup> /s)  | --              | <b>2.35E-06</b>          |
| Q/C <sub>vol</sub> | Inverse of mean concentration at the center of a 0.5 acre square source (g/m <sup>2</sup> -s per kg/m <sup>3</sup> ) | 68.18           | 68.18                    |
| T                  | Exposure interval (s)  | 9.5E+08         | 9.50E+08                 |
| ρ <sub>b</sub>     | Dry bulk soil density (g/cm <sup>3</sup> )   | 1.5             | 1.5                      |
| n                  | Total soil porosity 1-(ρ <sub>b</sub> /ρ <sub>s</sub> )  | 0.43            | 0.43                     |
| Θ <sub>a</sub>     | Air-filled soil porosity (n-Θ <sub>w</sub> )   | 0.17            | 0.17                     |
| Θ <sub>w</sub>     | Water-filled soil porosity   | 0.26            | 0.26                     |
| ρ <sub>s</sub>     | Soil particle density (g/cm <sup>3</sup> )   | 2.65            | 2.65                     |
| D <sub>a</sub>     | Diffusivity in air (cm <sup>2</sup> /s)  | --              | <b>0.06</b>              |
| H'                 | Dimensionless Henry's Law Constant   | --              | <b>0.03</b>              |
| D <sub>w</sub>     | Diffusivity in water (cm <sup>2</sup> /s)  | --              | <b>1.00E-05</b>          |
| K <sub>d</sub>     | Soil-water partition coefficient (cm <sup>3</sup> /g) = K <sub>oc</sub> × f <sub>oc</sub> (organics)                 | --              | <b>7.518</b>             |
| K <sub>oc</sub>    | Soil organic carbon partition coefficient (cm <sup>3</sup> /g)   | --              | <b>5012</b>              |
| f <sub>oc</sub>    | Fraction organic carbon in soil (g/g)  | 0.0015          | 0.0015                   |

Equations and Default values from New Mexico Environment Department's Technical Background Document for Development of Soil Screening Levels, Revisions 5.0, August 2009  
Physical and chemical properties from Massachusetts Department of Environmental Protection Massachusetts Contingency Plan Standards Spreadsheets 2009 (**bolded values**)

# Equation 16

## Derivation of the Particulate Emission Factor Residential and Commercial/Industrial Scenarios

$$PEF = Q / C_{wind} \times \frac{3,600 \text{ sec/hr}}{0.036 \times (1 - V) \times \left( \frac{U_m}{U_t} \right)^3 \times F(x)}$$

| Parameter | Definition (units)   | Default  |
|-----------|--|----------|
| PEF       | Particulate emission factor (m <sup>3</sup> /kg)   | 6.61E+09 |
| Q/Cwind   | Inverse of mean concentration at center of 0.5 acre square source (g/m <sup>2</sup> -sec per kg/m <sup>3</sup> ) | 81.85    |
| V         | Fraction of vegetative cover (unitless)  | 0.5      |
| Um        | Mean annual windspeed (m/s)  | 4.02     |
| Ut        | Equivalent threshold value of windspeed at 7 m (m/s)   | 11.32    |
| F(x)      | Function dependent on Um/Ut using Cowherd et al. 1985 (unitless)   | 0.0553   |

Equations and Default values from New Mexico Environment Department's Technical Background Document for Development of Soil Screening Levels, Revisions 5.0, August 2009







Western Refining - Bloomfield

**METALS ANALYSES**

| 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       |
| Cobalt    | SW-846 method 6010/6020       |
| Cyanide   | SW-846 method 335.3/335.2 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       |

# Chain-of-Custody Record

Client: Western Refining

Bloomfield Refined

Mailing Address: 50 Road 4990

Bloomfield NM 87413

Phone #: 505-632-4166

email or Fax#: Kelly.Robinson@WRCO

QA/QC Package: Level 4 (Full Validation)

☒ Standard ☐ Other

Accreditation: ☒ NELAP ☐ Other

Method (Type): EXCEL

Date: 8/15/11 Time: 940

Matrix: Soil

Sample Request ID: AOC 21-2 (0-0.5')

Container Type and #: 3 Jars

Preservative Type: None

TPH Method 8015B (Gas/Diesel)

BTEX + MTBE + TPH (Gas only)

TPH (Method 418.1)

EDB (Method 504.1)

8310 (PNA or PAH)

RCRA 8 Metals

Anions (F, Cl, NO<sub>3</sub>, NO<sub>2</sub>, PO<sub>4</sub>, SO<sub>4</sub>)

8081 Pesticides / 8082 PCBs

8260B (VOA), TPH - CRO

8270 (Semi-VOA)

TPH - MRO, DRO

Metals (See Method)

Air Bubbles (Y or N)

Turn-Around Time:

☒ Standard ☐ Rush

Project Name:

Group 6 RCAF LW,

Project #:

AOC No. 21

Project Manager:

Kelly Robinson

Sampler: Kelly & Tracy

On Site Temperature:

Container Type and #

Preservative Type

TPH Method 8015B (Gas/Diesel)

BTEX + MTBE + TPH (Gas only)

TPH (Method 418.1)

EDB (Method 504.1)

8310 (PNA or PAH)

RCRA 8 Metals

Anions (F, Cl, NO<sub>3</sub>, NO<sub>2</sub>, PO<sub>4</sub>, SO<sub>4</sub>)

8081 Pesticides / 8082 PCBs

8260B (VOA), TPH - CRO

8270 (Semi-VOA)

TPH - MRO, DRO

Metals (See Method)

Air Bubbles (Y or N)

Remarks:

See attached for Metals List

Received by: TD UPS

Date: 8/15/11 Time: 10:32



**HALL ENVIRONMENTAL  
ANALYSIS LABORATORY**

www.hallenviro.com

4901 Hawkins NE - Albuquerque, NM 87109

Tel. 505-345-3975 Fax 505-345-4107

Analysis Request

If necessary, samples submitted to Hall Environmental may be subcontracted to other accredited laboratories. Hall Environmental assumes as notice of this possibility. Any sub-contracted data will be clearly notated on the analytical report.

Western Refining - Bloomfield

**METALS ANALYSES**

| 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       |
| Cobalt    | SW-846 method 6010/6020       |
| Cyanide   | SW-846 method 335.3/335.2 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       |









# METALS ANALYSES

| 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       |
| Cobalt    | SW-846 method 6010/6020       |
| Cyanide   | SW-846 method 335.3/335.2 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       |

# Chain-of-Custody Record

Client: Western Refining  
Bloomfield, Refinery  
 Mailing Address: 50 Road 4990  
Bloomfield, NM 87413  
 Phone #: 505-632-4166  
 email or Fax: Kelly Robinson  
 QA/QC Package:  
☐ Standard ☒ Level 4 (Full Validation)  
 Accreditation  
☐ NELAP ☐ Other  
 XEDD (Type) EXCEL

| Date    | Time | Matrix | Sample Request ID      | Container Type and # | Preservative Type |
|---------|------|--------|------------------------|----------------------|-------------------|
| 8/16/11 | 0915 | Soil   | APC 19-17 (0-0.5')     | 3 Jan                | None              |
|         |      |        |                        | 2 empty              |                   |
|         |      |        |                        | 2 vials              | MeOH              |
|         |      | Soil   | APC 19-17 (0-0.5') DVA | 3 Jan                | None              |
|         |      |        |                        | 2 empty              |                   |
|         |      |        |                        | 2 vials              | MeOH              |
| 0945    |      | Soil   | APC 19-23 (0-0.5')     | 3 Jan                | None              |
|         |      |        |                        | 2 empty              |                   |
|         |      |        |                        | 2 vials              | MeOH              |
|         |      | Soil   | APC 19-21 (0-0.5')     | 3 Jan                | None              |
|         |      |        |                        | 2 empty              |                   |
|         |      |        |                        | 2 vials              | MeOH              |

Date: 8/16/11 Time: 1400  
 Relinquished by: Kelly Robinson  
 Date: 8/16/11 Time: 1400  
 Relinquished by: Kelly Robinson

Turn-Around Time:

☒ Standard ☐ Rush

Project Name:

GROUP 6 REA-LMV

Project #:

Project Manager:

Kelly Robinson

Sampler: Kelly Robinson

| BTEX + MTBE + TMBs (8021) | BTEX + MTBE + TPH (Gas only) | TPH Method 8015B (Gas/Diesel) | TPH (Method 418.1) | EDB (Method 504.1) | 8310 (PNA or PAH) | RCRA 8 Metals | Anions (F, Cl, NO <sub>3</sub> , NO <sub>2</sub> , PO <sub>4</sub> , SO <sub>4</sub> ) | 8081 Pesticides / 8082 PCBs | 8260B (VOA), TPH - GED | 8270 (Semi-VOA) | Metals (see attached) | TPH - GED, DRO, MRO | Air Bubbles (Y or N) |
|---------------------------|------------------------------|-------------------------------|--------------------|--------------------|-------------------|---------------|--|-----------------------------|------------------------|-----------------|-----------------------|---------------------|----------------------|
| -1                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -1                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -1                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -2                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -2                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -2                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -3                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -3                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -3                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -4                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -4                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |
| -4                        |                              |                               |                    |                    |                   |               |  |                             | X                      | X               | X                     | X                   |                      |

Remarks:

See attached for metals list

If necessary, samples submitted to Hall Environmental may be subcontracted to other accredited laboratories. This service as notice of this possibility. Any sub-contracted data will be clearly notated on the analytical report.

# Chain-of-Custody Record

Client: Western Refining  
Bloomfield Refinery  
 Mailing Address: 50 Road 4990  
Bloomfield, NM 87413  
 Phone #: 505-1632-4166  
 email or Fax: Kelly, Robinson@wlr.com  
 QA/QC Package:  
☐ Standard ☒ Level 4 (Full Validation)  
 Accreditation  
☐ NELAP ☐ Other \_\_\_\_\_  
 EDD (Type) F-Xcel

| Date    | Time | Matrix | Sample Request ID  | Container Type and # | Preservative Type |
|---------|------|--------|--------------------|----------------------|-------------------|
| 8/16/11 | 1025 | Soil   | APC 19-22 (0-0.5") | 3 Jar                | None              |
|         |      | ↓      | ↓                  | 2 evap               | ↓                 |
|         |      | ↓      | ↓                  | 2 vial               | MeOH              |
| 1045    |      | Soil   | APC 19-20 (0-0.5") | 3 Jar                | None              |
|         |      | ↓      | ↓                  | 2 evap               | ↓                 |
|         |      | ↓      | ↓                  | 2 vial               | MeOH              |
| 1105    |      | Soil   | APC 19-19 (0-0.5") | 3 Jar                | None              |
|         |      | ↓      | ↓                  | 2 evap               | ↓                 |
|         |      | ↓      | ↓                  | 2 vial               | MeOH              |
| ---     |      | ---    | MeOH Blank         | 2 vial               | MeOH              |

Date: 8/16/11 Time: 1400  
 Relinquished by: Kelly Robinson  
 Date: \_\_\_\_\_ Time: \_\_\_\_\_  
 Relinquished by: \_\_\_\_\_

Turn-Around Time: \_\_\_\_\_  
☒ Standard ☐ Rush  
 Project Name: Group 6 RRA 1W,  
 Project #: \_\_\_\_\_

Project Manager: Kelly Robinson  
 Sampler: Kelly Tracy

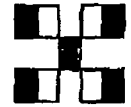
| BTEX + MTBE + TMBs (8021) | BTEX + MTBE + TPH (Gas only) | TPH Method 8015B (Gas/Diesel) | TPH (Method 418.1) | EDB (Method 504.1) | 8310 (PNA or PAH) | RCRA 8 Metals | Anions (F, Cl, NO <sub>3</sub> , NO <sub>2</sub> , PO <sub>4</sub> , SO <sub>4</sub> ) | 8081 Pesticides / 8082 PCB's | 8270 (Semi-VOA) | TPH-DPO, MRO | Metals (See attached) | Air Bubbles (Y or N) |
|---------------------------|------------------------------|-------------------------------|--------------------|--------------------|-------------------|---------------|--|------------------------------|-----------------|--------------|-----------------------|----------------------|
| -5                        |                              |                               |                    |                    |                   |               |  | X                            | X               | X            | X                     |                      |
| -5                        |                              |                               |                    |                    |                   |               |  | X                            | X               | X            | X                     |                      |
| -5                        |                              |                               |                    |                    |                   |               |  | X                            | X               | X            | X                     |                      |
| -6                        |                              |                               |                    |                    |                   |               |  | X                            | X               | X            | X                     |                      |
| -6                        |                              |                               |                    |                    |                   |               |  | X                            | X               | X            | X                     |                      |
| -7                        |                              |                               |                    |                    |                   |               |  | X                            | X               | X            | X                     |                      |
| -7                        |                              |                               |                    |                    |                   |               |  | X                            | X               | X            | X                     |                      |
| -8                        |                              |                               |                    |                    |                   |               |  | X                            | X               | X            | X                     |                      |

Received by: TO UPS  
 Date: 8/17/11 Time: 945  
 Received by: \_\_\_\_\_  
 Date: \_\_\_\_\_ Time: \_\_\_\_\_

Remarks: See Attached for Metals List

\* Necessary, samples submitted to Hall Environmental may be subcontracted to other accredited laboratories. This serves as notice of this possibility. Any sub-contracted data will be clearly noted on the analytical report.

292



**HALL ENVIRONMENTAL ANALYSIS LABORATORY**

www.hallenvironmental.com

4901 Hawkins NE - Albuquerque, NM 87109

Tel. 505-345-3975 Fax 505-345-4107

Analysis Request

Order #3

# METALS ANALYSES

| 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       |
| Cobalt    | SW-846 method 6010/6020       |
| Cyanide   | SW-846 method 335.3/335.2 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       |



## METALS ANALYSES

| 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        |
| Cobalt    | SW-846 method 6010/6020        |
| Cyanide   | SW-846 method 335.3/336.2 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        |

|   |   |                               |  |
|---|---|-------------------------------|--|
| <b>Chain-of-Custody Record</b>              |   | Turn-Around Time:             |  |
| Client: <u>Western Refining</u>             | <input checked="" type="checkbox"/> Standard                  | <input type="checkbox"/> Rush |  |
| <u>Bloomfield Refining</u>                  | Project Name:   |                               |  |
| Mailing Address: <u>50 Road 4990</u>        | <u>Group 6 RCRA I.W.</u>                                      |                               |  |
| <u>Bloomfield, NM 87413</u>                 | Project #:  |                               |  |
| Phone #: <u>505-632-4166</u>                | <u>AOC No. 19</u>   |                               |  |
| email or Fax: <u>Kelly.Robinson@wnr.com</u> | Project Manager:  |                               |  |
| QA/QC Package:                              | <u>Kelly Robinson</u>   |                               |  |
| <input type="checkbox"/> Standard           | <input checked="" type="checkbox"/> Level 4 (Full Validation) |                               |  |

### **Tum-Around Time:**

☒ Standard ☐ Rush

Project Name:

Group 6 RCEA LW.

Project #: AOC No. 19

Project Manager *Shelley Robinson*

Sampler: Kelly; Tracy

| Date | Time | Matrix | Sample Request ID |
|------|------|--------|-------------------|
|------|------|--------|-------------------|

## Sample Request ID

## Matrix

Date \_\_\_\_\_ Time \_\_\_\_\_

|         |      |      |                   |
|---------|------|------|-------------------|
| 8/16/11 | 1300 | Soil | Dec 19-7 (0-0.5') |
|---------|------|------|-------------------|

[illegible]

|         |      |      |                     |
|---------|------|------|---------------------|
| 8/16/11 | 1305 | Soil | POC 19-7 (4.5-2.0') |
|---------|------|------|---------------------|

[illegible]

|           |                   |
|-----------|-------------------|
| 1320 Soil | Box 19-9 (0-0.5') |
|-----------|-------------------|

[illegible]

|      |                   |
|------|-------------------|
| 1345 | Aug 19-18 (0-0.5) |
|------|-------------------|

|       |         |       |      |                  |                |
|-------|---------|-------|------|------------------|----------------|
| Date: | 8/17/11 | Time: | 1400 | Relinquished by: | Kelly Robinson |
|-------|---------|-------|------|------------------|----------------|

|       |       |                  |
|-------|-------|------------------|
| Date: | Time: | Relinquished by: |
|-------|-------|------------------|

| Date | Time |
|------|------|
|------|------|

Received by:

| Date | Time |
|------|------|
|------|------|

Received by-

Remarks:

Remarks:  
See Attached for Notepad List.

If necessary, samples submitted to Hall Environmental may be subcontracted to other accredited laboratories. This serves as notice of this possibility. Any sub-contracted data will be clearly notated on the analytical report.





# Chain-of-Custody Record

Client: Western Refining  
Bloomfield Refinery  
 Mailing Address: 50 Road 4990  
Bloomfield, NM 87413  
 Phone #: 505-632-4166

email or Fax#: Kelly.Robinson@wnr.com  
 QA/QC Package:  
☐ Standard ☒ Level 4 (Full Validation)  
 Accreditation  
☐ NELAP ☐ Other

Project Name: Group 6 RCEA Inv.  
 Project #:                       
 Project Manager: Kelly Robinson  
 Sampler: Kelly + Tracy  
 On Site:                       
 Sample Temperature:                     

| Date    | Time | Matrix | Sample Request ID | Container Type and # | Preservative Type | Turn-Around Time: <input checked="" type="checkbox"/> Standard <input type="checkbox"/> Rush |
|---------|------|--------|-------------------|----------------------|-------------------|--|
| 8/17/11 | 1045 | Soil   | AOC 20-8(0-0.5)   | 3 bu                 | None              |  |
|         |      |        |                   | 2 euro               |                   |  |
|         |      |        |                   | 2 vials              | MeOH              |  |
|         | 1120 | Soil   | AOC 20-9(0-0.5)   | 3 bu                 | None              |  |
|         |      |        |                   | 2 euro               |                   |  |
|         |      |        |                   | 2 vials              | MeOH              |  |
|         | 1140 | Soil   | AOC 20-10(0-0.5)  | 3 Jan                | None              |  |
|         |      |        |                   | 2 euro               |                   |  |
|         |      |        |                   | 2 vials              | MeOH              |  |
|         | 1200 | Soil   | AOC 20-11(0-0.5)  | 3 Jan                | None              |  |
|         |      |        |                   | 2 euro               |                   |  |
|         |      |        |                   | 2 vials              | MeOH              |  |

Date: 8/17/11 Time: 1400  
 Relinquished by: Kelly Robinson  
 Date: 8/18/11 Time: 1000  
 Relinquished by:                     

Received by: TO WRS  
 Date: 8/18/11 Time: 1000  
 Received by:                     

122  
**HALL ENVIRONMENTAL ANALYSIS LABORATORY**  
 www.hallenvironmental.com  
 4801 Hawkins NE - Albuquerque, NM 87109  
 Tel. 505-345-3975 Fax. 505-345-4107  
 Cooler #6

| Analysis Request   | TPH Method 418.1 | EDB (Method 504.1) | 8310 (PNA or PAH) | RCRA 8 Metals | Anions (F, Cl, NO <sub>3</sub> , NO <sub>2</sub> , PO <sub>4</sub> , SO <sub>4</sub> ) | 8081 Pesticides / 8082 PCBs | 8260B (EPA), TPH - GPO | 8270 (Semi-VOA) | TPH (DRO, GPO, GMR) | X Metals (See Attached) | Air Bubbles (Y or N) |
|--|------------------|--------------------|-------------------|---------------|--|-----------------------------|------------------------|-----------------|---------------------|-------------------------|----------------------|
| BTEX + MTBE + TMBs (8021)  |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| BTEX + MTBE + TPH (Gas only)   |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| TPH Method 8015B (Gas/Diesel)  |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| TPH (Method 418.1)   |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| EDB (Method 504.1)   |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| 8310 (PNA or PAH)  |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| RCRA 8 Metals  |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| Anions (F, Cl, NO <sub>3</sub> , NO <sub>2</sub> , PO <sub>4</sub> , SO <sub>4</sub> ) |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| 8081 Pesticides / 8082 PCBs  |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| 8260B (EPA), TPH - GPO   |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| 8270 (Semi-VOA)  |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| TPH (DRO, GPO, GMR)  |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |
| X Metals (See Attached)  |                  |                    |                   |               |  |                             |                        |                 |                     |                         |                      |

Remarks: See Attached for Metals List

If necessary, samples submitted to Hall Environmental may be subcontracted to other accredited laboratories. This serves as notice of this possibility. Any sub-contracted data will be clearly notated on the analytical report.



# METALS ANALYSES

| 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       |
| Cobalt    | SW-846 method 6010/6020       |
| Cyanide   | SW-846 method 335.3/335.2 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       |

# Chain-of-Custody Record

|   |  |  |  |
|---|--|--|--|
| Client: <u>Western Refining</u>   |  | Turn-Around Time: <input checked="" type="checkbox"/> Standard <input type="checkbox"/> Rush |  |
| Mailing Address: <u>Bloomfield Refinery</u>   |  | Project Name: <u>Group 6 RCTA Inv.</u>   |  |
| Phone #: <u>505-632-4166</u>  |  | Project #: <u>                    </u>   |  |
| email of Fax#: <u>Kelly, Robinson@lor.com</u>   |  | Project Manager: <u>Kelly Robinson</u>   |  |
| QA/QC Package: <u>Level 4 (Full Validation)</u>   |  | Sample: <u>Kelly Tracy</u>   |  |
| <input type="checkbox"/> Standard <input type="checkbox"/> NELAP <input type="checkbox"/> Other |  | On location: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No             |  |
| ACREDITED (Type) <u>EXCEL</u>   |  | Sample Temperature: <u>73</u>  |  |

| Date    | Time | Matrix | Sample Request ID     | Container Type and # | Preservative Type | TPH Method 8015B (Gas/Diesel) | TPH (Method 418.1) | EDB (Method 504.1) | 8310 (PNA or PAH) | RCRA 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 (MAY), TPH-CRO | 8270 (Semi-VOA) | TPH-DRO (MAY) | Metals (See attached) | Air Bubbles (Y or N) |
|---------|------|--------|-----------------------|----------------------|-------------------|-------------------------------|--------------------|--------------------|-------------------|---------------|--|------------------------------|----------------------|-----------------|---------------|-----------------------|----------------------|
| 8/17/11 | 840  | Soil   | ROC 20-3 (0-0.5')     | 3 Jars               | None              |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         |      |        |                       | 2 vials              | ↓                 |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         | 840  |        | ROC 20-3 (0-0.5') DUP | 3 Jars               | MeOH              |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         |      |        |                       | 2 vials              | ↓                 |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         |      |        |                       | 2 vials              | MeOH              |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         | 0900 |        | ROC 20-4 (0-0.5')     | 3 Jars               | None              |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         |      |        |                       | 2 vials              | ↓                 |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         |      |        |                       | 2 vials              | MeOH              |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         | 0945 |        | ROC 20-5 (0-0.5')     | 3 Jars               | None              |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         |      |        |                       | 2 vials              | ↓                 |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |
|         |      |        |                       | 2 vials              | MeOH              |                               |                    |                    |                   |               |  |                              |                      |                 |               |                       |                      |

|         |       |                  |              |         |       |
|---------|-------|------------------|--------------|---------|-------|
| Date:   | Time: | Relinquished by: | Received by: | Date:   | Time: |
| 8/17/11 | 1400  | Kelly Robinson   | To VES       | 8/18/11 | 1000  |
| Date:   | Time: | Relinquished by: | Received by: | Date:   | Time: |

Remarks: See attached for metals list



### METALS ANALYSES

| 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       |
| Cobalt    | SW-846 method 6010/6020       |
| Cyanide   | SW-846 method 335.3/335.2 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       |







## **1.0 DATA VALIDATION INTRODUCTION**

This summary presents data verification results for soil samples collected from soil borings installed at the Bloomfield Refinery in accordance with the approved Investigation Work Plan - Group 6. The data review was performed in accordance with the procedures specified in the Order issued by NMED (NMED, 2007), 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 45 soil samples were collected in August 2011 in accordance with the Group 6 Investigation Work Plan. Soil 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 metals (Antimony, arsenic, barium, beryllium, cadmium, chromium, cobalt, lead, nickel, selenium, silver, vanadium, and zinc) by SW846 Method 6010B;
- Cyanide by EPA Method 335.4; and
- Mercury by EPA Method 7471.

Additionally, 15 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 sample identifications, laboratory sample identifications, and requested analytical parameters.

## **2.0 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 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.

### **2.1 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.

## **2.2 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.

## **2.3 HOLDING TIMES**

All samples were extracted and analyzed within method-specified holding time limits.

## **2.4 BLANK CONTAMINATION**

### **2.4.1 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:

- 2-Butanone was detected in method blank for analytical batch 28144. Associated field sample results were non-detect; therefore data qualification was not needed.

### **2.4.2 Trip Blank**

Trip blanks were analyzed at the appropriate frequency as specified in the Order. Target compounds were not detected in the trip blanks.

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

Field and equipment rinsate blanks were collected at the appropriate frequency as specified in the Group 6 Investigation Work Plan. Target compounds were not detected in the field blanks and equipment rinsate blank, with the exception of the following:

- Bromodichlormethane was detected in equipment blank EB-AOC19, EB-AOC20, and EB-AOC21 at 2.3 ug/L, 2.6 ug/L, and 1.9 ug/L, respectively. Associated field samples were non-detect; therefore data qualification was not needed.
- Dibromochloromethane was detected in equipment blank EB-AOC19, EB-AOC20, and EB-AOC21 at 2.5 ug/L, 3.7 ug/L, and 2.7 ug/L, respectively. Associated field samples were non-detect; therefore data qualification was not needed.
- 2-Butanone was detected in equipment blank EB-AOC19 at 24 ug/L. Associated field sample AOC 19-23 (0-0.5') was qualified "J+" due to a potential high bias.
- Acetone was detected in equipment blank EB-AOC19 at 35 ug/L. Associated field sample AOC 19-23 (0-0.5') was qualified "J+" due to a potential high bias.

#### **2.4.4 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. Concentrations were not detected in method blanks, and therefore no field data was qualified due to common laboratory impacts.

#### **2.4.5 Methanol Blanks**

Methanol Blanks provided by the laboratory were analyzed for VOCs.

### **2.5 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:

- Surrogate recoveries for Di-n-octylphthalate (DNOP) and bromofluorobenzene (BFB) were below the lower acceptance limit for several field soil samples. Low surrogate recovery was due to required sample dilution for analytical analysis and/or matrix interference; therefore data qualification was not required.
- Surrogate recovery for 4-Bromofluorobenzene (18.4%) was below the lower acceptance limit of 70.0% for sample AOC 19-1 (0-0.5'). Data qualification was not required because remaining acid and base/neutral fractions were within acceptance limits.
- Surrogate recovery for 1,2-Dichloroethane-d4 was below the lower acceptance limit of 70% for several samples. Data qualification was not required because remaining acid and base/neutral fractions were within acceptance limits.
- Surrogate recovery for 4-Terphenyl-d14 (117%) was above the upper acceptance limit of 108% for Sample AOC 19-23 (0-0.5'). Data qualification was not required because the remaining acid and base/neutral fractions were within acceptance limits.
- Surrogate recovery for 4-Bromofluorobenzene (66.5%) was below the lower acceptance limit of 70.0% for sample AOC 20-3 (0-0.5'). Data qualification was not required because remaining acid and base/neutral fractions were within acceptance limits.
- Surrogate recovery for 4-Bromofluorobenzene (60.8%) was below the lower acceptance limit of 70.0% for sample AOC 20-3 (0-0.5') DUP. Data qualification was not required because remaining acid and base/neutral fractions were within acceptance limits.
- Surrogate recovery for DNOP (137%) was above the upper acceptance limit of 73.4% for AOC 21-5 (0-0.5'). The associated field sample result for DRO was qualified "J+" for a potential high bias.

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

LCS/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.
- 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; therefore no data qualification due to LCS/LCSD was required.

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

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 inorganic parameters result in sample qualification of the associated analytical batch.
- 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:

- The MS/MSD percent recoveries for Antimony (38.4% / 37.8%) were below the lower acceptance limit of 75% for analytical batch 28094. Associated field sample results for Antimony were non-detect. Data qualification "UJ" was required to indicate a potential bias for the associated samples.

- The MS/MS duplicate percent recoveries for Antimony (55.4%/ 50.1%) were below the lower acceptance limit of 75% for analytical batch 28123. The following associated field samples were qualified "UJ" to indicate a potential low bias:

|                      |                          |
|----------------------|--------------------------|
| - AOC 19-3 (0-0.5')  | - AOC 19-17 (0-0.5')     |
| - AOC 19-1 (0-0.5')  | - AOC 19-17 (0-0.5') DUP |
| - AOC 19-2 (0-0.5')  | - AOC 19-23 (0-0.5')     |
| - AOC 19-4 (0-0.5')  | - AOC 19-21 (0-0.5')     |
| - AOC 19-5 (0-0.5')  | - AOC 19-22 (0-0.5')     |
| - AOC 19-6 (0-0.5')  | - AOC 19-20 (0-0.5')     |
| - AOC 19-19 (0-0.5') |                          |

- The MS/MS duplicate percent recoveries for Selenium (56.9% / 61.3%) were below the lower acceptance limit of 75% for analytical batch 28123. The following associated field samples were qualified "UJ" to indicate a potential low bias:

|                      |                          |
|----------------------|--------------------------|
| - AOC 19-19 (0-0.5') | - AOC 19-17 (0-0.5')     |
| - AOC 19-21 (0-0.5') | - AOC 19-17 (0-0.5') DUP |
| - AOC 19-22 (0-0.5') | - AOC 19-23 (0-0.5')     |
| - AOC 19-20 (0-0.5') |                          |

## 2.8 DUPLICATES

### 2.8.1 Field Duplicates

Field duplicates were collected at a rate of 10 percent and submitted for analysis. 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.

Field duplicate RPDs were within acceptance limits except for the following:

- Barium for field sample AOC 21-1 (0-0.5');
- Zinc for field sample AOC 19-17;
- Diesel Range Organics (DRO), Motor Oil Range Organics (MRO), carbon disulfide, and beryllium for field sample AOC 20-3 (0-0.5').

### 3.0 COMPLETENESS SUMMARY

---

Two types of completeness were calculated for this project: contract and technical. The following equations were used to calculate the two types of completeness:

$$\% \text{ Contract Completeness} = \left( \frac{\text{Number of contract compliant results}}{\text{Number of reported results}} \right) \times 100$$

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

The overall contract completeness, which includes the evaluation of protocol and contract deviations, which includes the evaluation of the QC parameters listed in Section 2.0, was approximately 95 percent for soil analysis. The technical completeness attained for Group 6 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 Group 6 Work Plan were considered usable for the intended purposes and the project DQOs have been met.

**TABLE A-1**  
**Sampling and Analysis Schedule**



**Table A-1**  
**Sampling and Analysis Schedule**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

| Sample ID              | Lab ID      | Date Collected | Sample Type |
|------------------------|-------------|----------------|-------------|
| AOC 19-3 (0-0.5')      | 1108623-01  | 8/15/2011      | N           |
| AOC 19-1 (0-0.5')      | 1108623-02  | 8/15/2011      | N           |
| AOC 19-2 (0-0.5')      | 1108623-03  | 8/15/2011      | N           |
| AOC 19-4 (0-0.5')      | 1108623-04  | 8/15/2011      | N           |
| AOC 19-5 (0-0.5')      | 1108623-05  | 8/15/2011      | N           |
| AOC 19-6 (0-0.5')      | 1108623-06  | 8/15/2011      | N           |
| MeOH Blank             | 1108623-07  | 8/15/2011      | MB          |
| AOC 21-2 (0-0.5')      | 1108624-01  | 8/15/2011      | N           |
| AOC 21-2 (0-0.5') DUP  | 1108624-02  | 8/15/2011      | FD          |
| AOC 21-3 (0-0.5')      | 1108624-03  | 8/15/2011      | N           |
| AOC 21-4 (0-0.5')      | 1108624-04  | 8/15/2011      | N           |
| AOC 19-8 (0-0.5')      | 1108694-01  | 8/15/2011      | N           |
| AOC 19-10 (0-0.5')     | 1108694-02  | 8/15/2011      | N           |
| AOC 19-11 (0-0.5')     | 1108694-03  | 8/15/2011      | N           |
| AOC 19-12 (0-0.5')     | 1108694-04  | 8/15/2011      | N           |
| AOC 19-13 (0-0.5')     | 1108694-05  | 8/15/2011      | N           |
| AOC 19-14 (0-0.5')     | 1108694-06  | 8/16/2011      | N           |
| AOC 19-15 (0-0.5')     | 1108694-07  | 8/16/2011      | N           |
| AOC 19-16 (0-0.5')     | 1108694-08  | 8/16/2011      | N           |
| MeOH Blank             | 1108694-09  | 8/16/2011      | MB          |
| AOC 19-17 (0-0.5')     | 1108698-01  | 8/16/2011      | N           |
| AOC 19-17 (0-0.5') DUP | 1108698-02  | 8/16/2011      | FD          |
| AOC 19-23 (0-0.5')     | 1108698-03  | 8/16/2011      | N           |
| AOC 19-21 (0-0.5')     | 1108698-04  | 8/16/2011      | N           |
| AOC 19-22 (0-0.5')     | 1108698-05  | 8/16/2011      | N           |
| AOC 19-20 (0-0.5')     | 1108698-06  | 8/16/2011      | N           |
| AOC 19-19 (0-0.5')     | 1108698-07  | 8/16/2011      | N           |
| Methanol Blank         | 1108698-08  | 8/16/2011      | MB          |
| AOC 20-1 (0-0.5')      | 1108747-01  | 8/17/2011      | N           |
| AOC 20-2 (0-0.5')      | 1108747-02  | 8/17/2011      | N           |
| AOC 19-7 (0-0.5')      | 1108748-01  | 8/16/2011      | N           |
| AOC 19-7 (1.5-2.0)     | 1108748-02  | 8/16/2011      | N           |
| AOC 19-9 (0-0.5)       | 1108748-03  | 8/16/2011      | N           |
| AOC 19-18 (0-0.5)      | 1108748-04  | 8/16/2011      | N           |
| EB-AOC 19              | 1108748-05  | 8/16/2011      | EB          |
| MEOH BLANK             | 1108748-06  | 8/16/2011      | MB          |
| TRIP BLANK             | 1108748-07  | 8/16/2011      | TB          |
| AOC 20-8 (0-0.5')      | 1108750-01  | 8/17/2011      | N           |
| AOC 20-9 (0-0.5')      | 1108750-02  | 8/17/2011      | N           |
| AOC 20-10 (0-0.5')     | 1108750-03  | 8/17/2011      | N           |
| AOC 20-11 (0-0.5')     | 1108750-04  | 8/17/2011      | N           |
| AOC 20-12 (0-0.5')     | 1108750-05  | 8/17/2011      | N           |
| EB-AOC 20              | 1108750-06  | 8/17/2011      | EB          |
| Methanol Blank         | 1108750-07  | 8/17/2011      | MB          |
| Trip Blank             | 1108750-08  | 8/17/2011      | TB          |
| AOC 20-3 (0-0.5')      | 1108790-01  | 8/17/2011      | N           |
| AOC 20-3 (0-0.5') DUP  | 1108790-02B | 8/17/2011      | FD          |
| AOC 20-4 (0-0.5')      | 1108790-03  | 8/17/2011      | N           |
| AOC 20-5 (0-0.5')      | 1108790-04  | 8/17/2011      | N           |
| AOC 20-6 (0-0.5')      | 1108790-05  | 8/17/2011      | N           |
| AOC 20-7 (0-0.5')      | 1108790-06  | 8/17/2011      | N           |
| Methanol Blank         | 1108790-07  | 8/17/2011      | MB          |
| AOC 21-1 (0-0.5')      | 1108822-01  | 8/17/2011      | N           |
| AOC 21-1 (1.5-2.0')    | 1108822-02  | 8/17/2011      | N           |
| AOC 21-1 (2.0-2.5')    | 1108822-03  | 8/17/2011      | N           |
| AOC 21-5 (0-0.5')      | 1108822-04  | 8/17/2011      | N           |
| AOC 21-5 (1.5-2.0')    | 1108822-05  | 8/17/2011      | N           |
| AOC 21-5 (2.0-2.5')    | 1108822-06  | 8/17/2011      | N           |
| EB-AOC21               | 1108822-07  | 8/17/2011      | EB          |
| MEOH BLANK             | 1108822-08  | 8/17/2011      | MB          |

**Notes:**

VOCs = Volatile Organic Compounds

N = Normal field sample

FD = Field duplicate

TB = Trip Blank

EB = Equipment Blank

MB = Methanol Blank

**TABLE A-2**  
**Qualified Data**

**Table A-2**  
**Qualified Data**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

| Sample ID             | Date Collected | Analyte                     | Result | Units     | Matrix | Qualifier | Comments                                    |
|-----------------------|----------------|-----------------------------|--------|-----------|--------|-----------|---|
| AOC 19-2 (0-0.5)      | 8/15/2011      | Antimony                    | < 12   | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-4 (0-0.5)      | 8/15/2011      | Antimony                    | < 13   | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-6 (0-0.5)      | 8/15/2011      | Antimony                    | < 13   | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-1 (0-0.5)      | 8/15/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-3 (0-0.5)      | 8/15/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-5 (0-0.5)      | 8/15/2011      | Antimony                    | < 5.0  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-3 (0-0.5)      | 8/16/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-1 (0-0.5)      | 8/16/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-2 (0-0.5)      | 8/16/2011      | Antimony                    | < 12   | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-4 (0-0.5)      | 8/16/2011      | Antimony                    | < 13   | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-5 (0-0.5)      | 8/16/2011      | Antimony                    | < 5.0  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-6 (0-0.5)      | 8/16/2011      | Antimony                    | < 13   | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-17 (0-0.5)     | 8/16/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-17 (0-0.5)     | 8/16/2011      | Selenium                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-17 (0-0.5) DUP | 8/16/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-17 (0-0.5) DUP | 8/16/2011      | Selenium                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-23 (0-0.5)     | 8/16/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-23 (0-0.5)     | 8/16/2011      | Selenium                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-21 (0-0.5)     | 8/16/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-21 (0-0.5)     | 8/16/2011      | Selenium                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-22 (0-0.5)     | 8/16/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-22 (0-0.5)     | 8/16/2011      | Selenium                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-20 (0-0.5)     | 8/16/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-20 (0-0.5)     | 8/16/2011      | Selenium                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-19 (0-0.5)     | 8/16/2011      | Antimony                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-19 (0-0.5)     | 8/16/2011      | Selenium                    | < 2.5  | mg/Kg     | Soil   | UJ        | Qualified due to low MS/ MSD recovery       |
| AOC 19-23 (0-0.5)     | 8/16/2011      | 2-Butanone                  | 43.8   | ug/kg-dry | Soil   | J+        | Qualified due to equipment blank detections |
| AOC 19-23 (0-0.5)     | 8/16/2011      | Acetone                     | 134    | ug/kg-dry | Soil   | J+        | Qualified due to equipment blank detections |
| AOC 21-5 (0-0.5)      | 8/17/2011      | Diesel Range Organics (DRO) | 57     | mg/kg     | Soil   | J+        | Qualified due to high surrogate recovery    |

MS/MSD - Matrix spike/matrix spike duplicate

ug/kg - microgram per kilogram

mg/kg - milligrams per kilogram

UJ - Estimated reporting limit

J+ = high bias

**TABLE A-3**  
**Field Duplicate Summary**

**Table A-3**  
**Field Duplicate Summary**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                  | Parameter                      | AOC 21-2 (0-0.5')<br>Sample Result | AOC 21-2 (0-0.5') DUP<br>Field Duplicate | RPD<br>(%) |
|------------------|--------------------------------|------------------------------------|--|------------|
| TPH (mg/kg-dry): | Diesel Range Organics (DRO)    | 710                                | 860                                      | 19.1       |
|                  | Motor Oil Range Organics (MRO) | < 510                              | < 500                                    | NC         |
|                  | Gasoline Range Organics (GRO)  | < 5.0                              | < 5.0                                    | NC         |
| VOCs (ug/kg-dry) | 1,1,1,2-Tetrachloroethane      | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1,1-Trichloroethane          | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1,2,2-Tetrachloroethane      | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1,2-Trichloroethane          | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1-Dichloroethane             | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1-Dichloroethene             | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1-Dichloropropene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2,3-Trichlorobenzene         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2,3-Trichloropropane         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2,4-Trichlorobenzene         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2,4-Trimethylbenzene         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2-Dibromo-3-chloropropane    | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2-Dibromoethane (EDB)        | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2-Dichlorobenzene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2-Dichloroethane (EDC)       | < 3.00                             | < 3.00                                   | NC         |
|                  | 1,2-Dichloropropane            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,3,5-Trimethylbenzene         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,3-Dichlorobenzene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,3-Dichloropropane            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,4-Dichlorobenzene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 2,2-Dichloropropane            | < 1.00                             | < 1.00                                   | NC         |
|                  | 2-Butanone                     | < 1.00                             | < 1.00                                   | NC         |
|                  | 2-Chlorotoluene                | < 1.00                             | < 1.00                                   | NC         |
|                  | 2-Hexanone                     | < 1.00                             | < 1.00                                   | NC         |
|                  | 4-Chlorotoluene                | < 1.00                             | < 1.00                                   | NC         |
|                  | 4-Isopropyltoluene             | < 1.00                             | < 1.00                                   | NC         |
|                  | 4-Methyl-2-pentanone           | < 1.00                             | < 1.00                                   | NC         |
|                  | Acetone                        | < 2.00                             | < 2.00                                   | NC         |
|                  | Benzene                        | < 1.00                             | < 1.00                                   | NC         |
|                  | Bromobenzene                   | < 1.00                             | < 1.00                                   | NC         |
|                  | Bromodichloromethane           | < 1.00                             | < 1.00                                   | NC         |
|                  | Bromoform                      | < 1.00                             | < 1.00                                   | NC         |
|                  | Bromomethane                   | < 1.00                             | < 1.00                                   | NC         |
|                  | Carbon disulfide               | < 1.00                             | < 1.00                                   | NC         |
|                  | Carbon tetrachloride           | < 1.00                             | < 1.00                                   | NC         |
|                  | Chlorobenzene                  | < 1.00                             | < 1.00                                   | NC         |
|                  | Chloroethane                   | < 1.00                             | < 1.00                                   | NC         |
|                  | Chloroform                     | < 1.00                             | < 1.00                                   | NC         |
|                  | Chloromethane                  | < 1.00                             | < 1.00                                   | NC         |
|                  | cis-1,2-DCE                    | < 1.00                             | < 1.00                                   | NC         |
|                  | cis-1,3-Dichloropropene        | < 1.00                             | < 1.00                                   | NC         |
|                  | Dibromochloromethane           | < 1.00                             | < 1.00                                   | NC         |
|                  | Dibromomethane                 | < 1.00                             | < 1.00                                   | NC         |
|                  | Dichlorodifluoromethane        | < 1.00                             | < 1.00                                   | NC         |
|                  | Ethylbenzene                   | < 1.00                             | < 1.00                                   | NC         |
|                  | Hexachlorobutadiene            | < 1.00                             | < 1.00                                   | NC         |
|                  | Isopropylbenzene               | < 1.00                             | < 1.00                                   | NC         |
|                  | Methyl tert-butyl ether (MTBE) | < 1.00                             | < 1.00                                   | NC         |
|                  | Methylene chloride             | < 1.00                             | < 1.00                                   | NC         |
|                  | Naphthalene                    | < 1.00                             | < 1.00                                   | NC         |
|                  | n-Butylbenzene                 | < 1.00                             | < 1.00                                   | NC         |
|                  | n-Propylbenzene                | < 1.00                             | < 1.00                                   | NC         |
|                  | sec-Butylbenzene               | < 1.00                             | < 1.00                                   | NC         |
|                  | Styrene                        | < 1.00                             | < 1.00                                   | NC         |
|                  | tert-Butylbenzene              | < 1.00                             | < 1.00                                   | NC         |
|                  | Tetrachloroethene (PCE)        | < 1.00                             | < 1.00                                   | NC         |
|                  | Toluene                        | 19.3                               | 25.6                                     | 28.1       |
|                  | trans-1,2-DCE                  | < 1.00                             | < 1.00                                   | NC         |
|                  | trans-1,3-Dichloropropene      | < 1.00                             | < 1.00                                   | NC         |
|                  | Trichloroethene (TCE)          | < 1.00                             | < 1.00                                   | NC         |
|                  | Trichlorofluoromethane         | < 1.00                             | < 1.00                                   | NC         |
|                  | Vinyl chloride                 | < 1.00                             | < 1.00                                   | NC         |
|                  | Xylenes, Total                 | < 1.00                             | < 1.00                                   | NC         |

**Table A-3**  
**Field Duplicate Summary**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                    | Parameter                   | AOC 21-2 (0-0.5')<br>Sample Result | AOC 21-2 (0-0.5') DUP<br>Field Duplicate | RPD<br>(%) |
|--------------------|-----------------------------|------------------------------------|--|------------|
| SVOCs (mg/kg-dry): | 1,2,4-Trichlorobenzene      | < 0.20                             | < 0.20                                   | NC         |
|                    | 1,2-Dichlorobenzene         | < 0.20                             | < 0.20                                   | NC         |
|                    | 1,3-Dichlorobenzene         | < 0.20                             | < 0.20                                   | NC         |
|                    | 1,4-Dichlorobenzene         | < 0.20                             | < 0.20                                   | NC         |
|                    | 1-Methylnaphthalene         | < 0.20                             | < 0.20                                   | NC         |
|                    | 2,4,5-Trichlorophenol       | < 0.20                             | < 0.20                                   | NC         |
|                    | 2,4,6-Trichlorophenol       | < 0.20                             | < 0.20                                   | NC         |
|                    | 2,4-Dichlorophenol          | < 0.40                             | < 0.40                                   | NC         |
|                    | 2,4-Dimethylphenol          | < 0.30                             | < 0.30                                   | NC         |
|                    | 2,4-Dinitrophenol           | < 0.40                             | < 0.40                                   | NC         |
|                    | 2,4-Dinitrotoluene          | < 0.50                             | < 0.50                                   | NC         |
|                    | 2,6-Dinitrotoluene          | < 0.50                             | < 0.50                                   | NC         |
|                    | 2-Chloronaphthalene         | < 0.25                             | < 0.25                                   | NC         |
|                    | 2-Chlorophenol              | < 0.20                             | < 0.20                                   | NC         |
|                    | 2-Methylnaphthalene         | < 0.20                             | < 0.20                                   | NC         |
|                    | 2-Methylphenol              | < 0.50                             | < 0.50                                   | NC         |
|                    | 2-Nitroaniline              | < 0.20                             | < 0.20                                   | NC         |
|                    | 2-Nitrophenol               | < 0.20                             | < 0.20                                   | NC         |
|                    | 3,3'-Dichlorobenzidine      | < 0.25                             | < 0.25                                   | NC         |
|                    | 3+4-Methylphenol            | < 0.20                             | < 0.20                                   | NC         |
|                    | 3-Nitroaniline              | < 0.20                             | < 0.20                                   | NC         |
|                    | 4,6-Dinitro-2-methylphenol  | < 0.50                             | < 0.50                                   | NC         |
|                    | 4-Bromophenyl phenyl ether  | < 0.20                             | < 0.20                                   | NC         |
|                    | 4-Chloro-3-methylphenol     | < 0.50                             | < 0.50                                   | NC         |
|                    | 4-Chloroaniline             | < 0.50                             | < 0.50                                   | NC         |
|                    | 4-Chlorophenyl phenyl ether | < 0.20                             | < 0.20                                   | NC         |
|                    | 4-Nitroaniline              | < 0.40                             | < 0.40                                   | NC         |
|                    | 4-Nitrophenol               | < 0.25                             | < 0.25                                   | NC         |
|                    | Acenaphthene                | < 0.20                             | < 0.20                                   | NC         |
|                    | Acenaphthylene              | < 0.20                             | < 0.20                                   | NC         |
|                    | Aniline                     | < 0.20                             | < 0.20                                   | NC         |
|                    | Anthracene                  | < 0.20                             | < 0.20                                   | NC         |
|                    | Azobenzene                  | < 0.20                             | < 0.20                                   | NC         |
|                    | Benz(a)anthracene           | < 0.20                             | < 0.20                                   | NC         |
|                    | Benzo(a)pyrene              | < 0.20                             | < 0.20                                   | NC         |
|                    | Benzo(b)fluoranthene        | < 0.20                             | < 0.20                                   | NC         |
|                    | Benzo(g,h,i)perylene        | < 0.20                             | < 0.20                                   | NC         |
|                    | Benzo(k)fluoranthene        | < 0.20                             | < 0.20                                   | NC         |
|                    | Benzoic acid                | < 0.50                             | < 0.50                                   | NC         |
|                    | Benzyl alcohol              | < 0.20                             | < 0.20                                   | NC         |
|                    | Bis(2-chloroethoxy)methane  | < 0.20                             | < 0.20                                   | NC         |
|                    | Bis(2-chloroethyl)ether     | < 0.20                             | < 0.20                                   | NC         |
|                    | Bis(2-chloroisopropyl)ether | < 0.20                             | < 0.20                                   | NC         |
|                    | Bis(2-ethylhexyl)phthalate  | < 0.50                             | < 0.50                                   | NC         |
|                    | Butyl benzyl phthalate      | < 0.20                             | < 0.20                                   | NC         |
|                    | Carbazole                   | < 0.20                             | < 0.20                                   | NC         |
|                    | Chrysene                    | < 0.20                             | < 0.20                                   | NC         |
|                    | Dibenz(a,h)anthracene       | < 0.20                             | < 0.20                                   | NC         |
|                    | Dibenzofuran                | < 0.20                             | < 0.20                                   | NC         |
|                    | Diethyl phthalate           | < 0.20                             | < 0.20                                   | NC         |
|                    | Dimethyl phthalate          | < 0.20                             | < 0.20                                   | NC         |
|                    | Di-n-butyl phthalate        | < 0.50                             | < 0.50                                   | NC         |
|                    | Di-n-octyl phthalate        | < 0.25                             | < 0.25                                   | NC         |
|                    | Fluoranthene                | < 0.20                             | < 0.20                                   | NC         |
|                    | Fluorene                    | < 0.20                             | < 0.20                                   | NC         |
|                    | Hexachlorobenzene           | < 0.20                             | < 0.20                                   | NC         |
|                    | Hexachlorobutadiene         | < 0.20                             | < 0.20                                   | NC         |
|                    | Hexachlorocyclopentadiene   | < 0.20                             | < 0.20                                   | NC         |
|                    | Hexachloroethane            | < 0.20                             | < 0.20                                   | NC         |
|                    | Indeno(1,2,3-cd)pyrene      | < 0.20                             | < 0.20                                   | NC         |
|                    | Isophorone                  | < 0.50                             | < 0.50                                   | NC         |
|                    | Naphthalene                 | < 0.20                             | < 0.20                                   | NC         |
|                    | Nitrobenzene                | < 0.50                             | < 0.50                                   | NC         |
|                    | N-Nitrosodi-n-propylamine   | < 0.20                             | < 0.20                                   | NC         |
|                    | N-Nitrosodiphenylamine      | < 0.20                             | < 0.20                                   | NC         |
|                    | Pentachlorophenol           | < 0.40                             | < 0.40                                   | NC         |
|                    | Phenanthrene                | < 0.20                             | < 0.20                                   | NC         |
|                    | Phenol                      | < 0.20                             | < 0.20                                   | NC         |
|                    | Pyrene                      | < 0.20                             | < 0.20                                   | NC         |
|                    | Pyridine                    | < 0.50                             | < 0.50                                   | NC         |

**Table A-3**  
**Field Duplicate Summary**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                     | Parameter | AOC 21-2 (0-0.5')<br>Sample Result | AOC 21-2 (0-0.5') DUP<br>Field Duplicate | RPD<br>(%) |
|---------------------|-----------|------------------------------------|--|------------|
| Metals (mg/kg-dry): | Antimony  | < 13                               | < 12                                     | NC         |
|                     | Arsenic   | < 13                               | < 12                                     | NC         |
|                     | Barium    | 28                                 | 45                                       | 46.6       |
|                     | Beryllium | 2.1                                | 2.1                                      | 0.0        |
|                     | Cadmium   | < 0.50                             | < 0.50                                   | NC         |
|                     | Chromium  | 11                                 | 11                                       | 0          |
|                     | Cobalt    | 8.5                                | 7.7                                      | 9.9        |
|                     | Cyanide   | < 0.30                             | < 0.30                                   | NC         |
|                     | Lead      | 10                                 | 11                                       | 9.5        |
|                     | Mercury   | < 0.033                            | < 0.033                                  | NC         |
|                     | Nickel    | 10                                 | 10                                       | 0.0        |
|                     | Selenium  | < 13                               | < 12                                     | NC         |
|                     | Silver    | < 1.3                              | < 1.2                                    | NC         |
|                     | Vanadium  | 17                                 | 16                                       | 6          |
|                     | Zinc      | 69                                 | 64                                       | 7.5        |

**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

\* = Field Duplicate RPD Outlier

**Table A-3**  
**Field Duplicate Summary**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                  | Parameter                      | AOC 19-17 (0-0.5')<br>Sample Result | AOC 19-17 (0-0.5') DUP<br>Field Duplicate | RPD<br>(%) |
|------------------|--------------------------------|-------------------------------------|---|------------|
| TPH (mg/kg-dry): | Diesel Range Organics (DRO)    | 19                                  | 21  | 10.0       |
|                  | Motor Oil Range Organics (MRO) | 56                                  | 63  | 11.8       |
|                  | Gasoline Range Organics (GRO)  | < 5.0                               | < 5.0                                     | NC         |
| VOCs (ug/kg-dry) | 1,1,1,2-Tetrachloroethane      | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,1,1-Trichloroethane          | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,1,2,2-Tetrachloroethane      | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,1,2-Trichloroethane          | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,1-Dichloroethane             | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,1-Dichloroethene             | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,1-Dichloropropene            | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,2,3-Trichlorobenzene         | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,2,3-Trichloropropane         | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,2,4-Trichlorobenzene         | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,2,4-Trimethylbenzene         | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,2-Dibromo-3-chloropropane    | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,2-Dibromoethane (EDB)        | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,2-Dichlorobenzene            | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,2-Dichloroethane (EDC)       | < 3.00                              | < 3.00                                    | NC         |
|                  | 1,2-Dichloropropane            | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,3,5-Trimethylbenzene         | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,3-Dichlorobenzene            | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,3-Dichloropropane            | < 1.00                              | < 1.00                                    | NC         |
|                  | 1,4-Dichlorobenzene            | < 1.00                              | < 1.00                                    | NC         |
|                  | 1-Methylnaphthalene            | < 1.00                              | < 1.00                                    | NC         |
|                  | 2,2-Dichloropropane            | < 1.00                              | < 1.00                                    | NC         |
|                  | 2-Butanone                     | < 1.00                              | < 1.00                                    | NC         |
|                  | 2-Chlorotoluene                | < 1.00                              | < 1.00                                    | NC         |
|                  | 2-Hexanone                     | < 1.00                              | < 1.00                                    | NC         |
|                  | 2-Methylnaphthalene            | < 1.00                              | < 1.00                                    | NC         |
|                  | 4-Chlorotoluene                | < 1.00                              | < 1.00                                    | NC         |
|                  | 4-Isopropyltoluene             | < 1.00                              | < 1.00                                    | NC         |
|                  | 4-Methyl-2-pentanone           | < 1.00                              | < 1.00                                    | NC         |
|                  | Acetone                        | < 2.00                              | < 2.00                                    | NC         |
|                  | Benzene                        | < 1.00                              | < 1.00                                    | NC         |
|                  | Bromobenzene                   | < 1.00                              | < 1.00                                    | NC         |
|                  | Bromodichloromethane           | < 1.00                              | < 1.00                                    | NC         |
|                  | Bromoform                      | < 1.00                              | < 1.00                                    | NC         |
|                  | Bromomethane                   | < 1.00                              | < 1.00                                    | NC         |
|                  | Carbon disulfide               | < 1.00                              | < 1.00                                    | NC         |
|                  | Carbon tetrachloride           | < 1.00                              | < 1.00                                    | NC         |
|                  | Chlorobenzene                  | < 1.00                              | < 1.00                                    | NC         |
|                  | Chloroethane                   | < 1.00                              | < 1.00                                    | NC         |
|                  | Chloroform                     | < 1.00                              | < 1.00                                    | NC         |
|                  | Chloromethane                  | < 1.00                              | < 1.00                                    | NC         |
|                  | cis-1,2-DCE                    | < 1.00                              | < 1.00                                    | NC         |
|                  | cis-1,3-Dichloropropene        | < 1.00                              | < 1.00                                    | NC         |
|                  | Dibromochloromethane           | < 1.00                              | < 1.00                                    | NC         |
|                  | Dibromomethane                 | < 1.00                              | < 1.00                                    | NC         |
|                  | Dichlorodifluoromethane        | < 1.00                              | < 1.00                                    | NC         |
|                  | Ethylbenzene                   | < 1.00                              | < 1.00                                    | NC         |
|                  | Hexachlorobutadiene            | < 1.00                              | < 1.00                                    | NC         |
|                  | Isopropylbenzene               | < 1.00                              | < 1.00                                    | NC         |
|                  | Methyl tert-butyl ether (MTBE) | < 1.00                              | < 1.00                                    | NC         |
|                  | Methylene chloride             | < 1.00                              | < 1.00                                    | NC         |
|                  | Naphthalene                    | < 1.00                              | < 1.00                                    | NC         |
|                  | n-Butylbenzene                 | < 1.00                              | < 1.00                                    | NC         |
|                  | n-Propylbenzene                | < 1.00                              | < 1.00                                    | NC         |
|                  | sec-Butylbenzene               | < 1.00                              | < 1.00                                    | NC         |
|                  | Styrene                        | < 1.00                              | < 1.00                                    | NC         |
|                  | tert-Butylbenzene              | < 1.00                              | < 1.00                                    | NC         |
|                  | Tetrachloroethene (PCE)        | < 1.00                              | < 1.00                                    | NC         |
|                  | Toluene                        | 4.42                                | 4.02                                      | 9.5        |
|                  | trans-1,2-DCE                  | < 1.00                              | < 1.00                                    | NC         |
|                  | trans-1,3-Dichloropropene      | < 1.00                              | < 1.00                                    | NC         |
|                  | Trichloroethene (TCE)          | < 1.00                              | < 1.00                                    | NC         |
|                  | Trichlorofluoromethane         | < 1.00                              | < 1.00                                    | NC         |
|                  | Vinyl chloride                 | < 1.00                              | < 1.00                                    | NC         |
|                  | Xylenes, Total                 | < 1.00                              | < 1.00                                    | NC         |



**Table A-3**  
**Field Duplicate Summary**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                    | Parameter                   | AOC 19-17 (0-0.5')<br>Sample Result | AOC 19-17 (0-0.5') DUP<br>Field Duplicate | RPD<br>(%) |
|--------------------|-----------------------------|-------------------------------------|---|------------|
| SVOCs (mg/kg-dry): | 1,2,4-Trichlorobenzene      | < 0.20                              | < 0.20                                    | NC         |
|                    | 1,2-Dichlorobenzene         | < 0.20                              | < 0.20                                    | NC         |
|                    | 1,3-Dichlorobenzene         | < 0.20                              | < 0.20                                    | NC         |
|                    | 1,4-Dichlorobenzene         | < 0.20                              | < 0.20                                    | NC         |
|                    | 1-Methylnaphthalene         | < 0.20                              | < 0.20                                    | NC         |
|                    | 2,4,5-Trichlorophenol       | < 0.20                              | < 0.20                                    | NC         |
|                    | 2,4,6-Trichlorophenol       | < 0.20                              | < 0.20                                    | NC         |
|                    | 2,4-Dichlorophenol          | < 0.40                              | < 0.40                                    | NC         |
|                    | 2,4-Dimethylphenol          | < 0.30                              | < 0.30                                    | NC         |
|                    | 2,4-Dinitrophenol           | < 0.40                              | < 0.40                                    | NC         |
|                    | 2,4-Dinitrotoluene          | < 0.50                              | < 0.50                                    | NC         |
|                    | 2,6-Dinitrotoluene          | < 0.50                              | < 0.50                                    | NC         |
|                    | 2-Chloronaphthalene         | < 0.25                              | < 0.25                                    | NC         |
|                    | 2-Chlorophenol              | < 0.20                              | < 0.20                                    | NC         |
|                    | 2-Methylnaphthalene         | < 0.20                              | < 0.20                                    | NC         |
|                    | 2-Methylphenol              | < 0.50                              | < 0.50                                    | NC         |
|                    | 2-Nitroaniline              | < 0.20                              | < 0.20                                    | NC         |
|                    | 2-Nitrophenol               | < 0.20                              | < 0.20                                    | NC         |
|                    | 3,3'-Dichlorobenzidine      | < 0.25                              | < 0.25                                    | NC         |
|                    | 3+4-Methylphenol            | < 0.20                              | < 0.20                                    | NC         |
|                    | 3-Nitroaniline              | < 0.20                              | < 0.20                                    | NC         |
|                    | 4,6-Dinitro-2-methylphenol  | < 0.50                              | < 0.50                                    | NC         |
|                    | 4-Bromophenyl phenyl ether  | < 0.20                              | < 0.20                                    | NC         |
|                    | 4-Chloro-3-methylphenol     | < 0.50                              | < 0.50                                    | NC         |
|                    | 4-Chloroaniline             | < 0.50                              | < 0.50                                    | NC         |
|                    | 4-Chlorophenyl phenyl ether | < 0.20                              | < 0.20                                    | NC         |
|                    | 4-Nitroaniline              | < 0.40                              | < 0.40                                    | NC         |
|                    | 4-Nitrophenol               | < 0.25                              | < 0.25                                    | NC         |
|                    | Acenaphthene                | < 0.20                              | < 0.20                                    | NC         |
|                    | Acenaphthylene              | < 0.20                              | < 0.20                                    | NC         |
|                    | Aniline                     | < 0.20                              | < 0.20                                    | NC         |
|                    | Anthracene                  | < 0.20                              | < 0.20                                    | NC         |
|                    | Azobenzene                  | < 0.20                              | < 0.20                                    | NC         |
|                    | Benzo(a)anthracene          | < 0.20                              | < 0.20                                    | NC         |
|                    | Benzo(a)pyrene              | < 0.20                              | < 0.20                                    | NC         |
|                    | Benzo(b)fluoranthene        | < 0.20                              | < 0.20                                    | NC         |
|                    | Benzo(g,h,i)perylene        | < 0.20                              | < 0.20                                    | NC         |
|                    | Benzo(k)fluoranthene        | < 0.20                              | < 0.20                                    | NC         |
|                    | Benzoic acid                | < 0.50                              | < 0.50                                    | NC         |
|                    | Benzyl alcohol              | < 0.20                              | < 0.20                                    | NC         |
|                    | Bis(2-chloroethoxy)methane  | < 0.20                              | < 0.20                                    | NC         |
|                    | Bis(2-chloroethyl)ether     | < 0.20                              | < 0.20                                    | NC         |
|                    | Bis(2-chloroisopropyl)ether | < 0.20                              | < 0.20                                    | NC         |
|                    | Bis(2-ethylhexyl)phthalate  | < 0.50                              | < 0.50                                    | NC         |
|                    | Butyl benzyl phthalate      | < 0.20                              | < 0.20                                    | NC         |
|                    | Carbazole                   | < 0.20                              | < 0.20                                    | NC         |
|                    | Chrysene                    | < 0.20                              | < 0.20                                    | NC         |
|                    | Dibenz(a,h)anthracene       | < 0.20                              | < 0.20                                    | NC         |
|                    | Dibenzofuran                | < 0.20                              | < 0.20                                    | NC         |
|                    | Diethyl phthalate           | < 0.20                              | < 0.20                                    | NC         |
|                    | Dimethyl phthalate          | < 0.20                              | < 0.20                                    | NC         |
|                    | Di-n-butyl phthalate        | < 0.50                              | < 0.50                                    | NC         |
|                    | Di-n-octyl phthalate        | < 0.25                              | < 0.25                                    | NC         |
|                    | Fluoranthene                | < 0.20                              | < 0.20                                    | NC         |
|                    | Fluorene                    | < 0.20                              | < 0.20                                    | NC         |
|                    | Hexachlorobenzene           | < 0.20                              | < 0.20                                    | NC         |
|                    | Hexachlorobutadiene         | < 0.20                              | < 0.20                                    | NC         |
|                    | Hexachlorocyclopentadiene   | < 0.20                              | < 0.20                                    | NC         |
|                    | Hexachloroethane            | < 0.20                              | < 0.20                                    | NC         |
|                    | Indeno(1,2,3-cd)pyrene      | < 0.20                              | < 0.20                                    | NC         |
|                    | Isophorone                  | < 0.50                              | < 0.50                                    | NC         |
|                    | Naphthalene                 | < 0.20                              | < 0.20                                    | NC         |
|                    | Nitrobenzene                | < 0.50                              | < 0.50                                    | NC         |
|                    | N-Nitrosodi-n-propylamine   | < 0.20                              | < 0.20                                    | NC         |
|                    | N-Nitrosodiphenylamine      | < 0.20                              | < 0.20                                    | NC         |
|                    | Pentachlorophenol           | < 0.40                              | < 0.40                                    | NC         |
|                    | Phenanthrene                | < 0.20                              | < 0.20                                    | NC         |
|                    | Phenol                      | < 0.20                              | < 0.20                                    | NC         |
|                    | Pyrene                      | < 0.20                              | < 0.20                                    | NC         |

**Table A-3**  
**Field Duplicate Summary**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                            | Parameter | AOC 19-17 (0-0.5')<br>Sample Result | AOC 19-17 (0-0.5') DUP<br>Field Duplicate | RPD<br>(%) |
|----------------------------|-----------|-------------------------------------|---|------------|
| <b>Metals (mg/kg-dry):</b> | Pyridine  | < 0.50                              | < 0.50                                    | NC         |
|                            | Antimony  | < 2.5                               | < 2.5                                     | NC         |
|                            | Arsenic   | 3.8                                 | < 2.5                                     | NC         |
|                            | Barium    | 120                                 | 120                                       | 0.0        |
|                            | Beryllium | 0.19                                | 0.25                                      | 27.3       |
|                            | Cadmium   | < 0.10                              | < 0.10                                    | NC         |
|                            | Chromium  | 33                                  | 43  | 26.3       |
|                            | Cobalt    | 1.4                                 | 1.8                                       | 25.0       |
|                            | Cyanide   | < 0.30                              | < 0.30                                    | NC         |
|                            | Lead      | 3.1                                 | 3.6                                       | 14.9       |
|                            | Mercury   | < 0.033                             | < 0.033                                   | NC         |
|                            | Nickel    | 1.6                                 | 2.1                                       | 27.0       |
|                            | Selenium  | < 2.5                               | < 2.5                                     | NC         |
|                            | Silver    | < 0.25                              | < 0.25                                    | NC         |
|                            | Vanadium  | 7.2                                 | 8.8                                       | 20.0       |
|                            | Zinc      | 43                                  | 77  | 56.7       |

**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

\* = Field Duplicate RPD Outlier

**Table A-3**  
**Field Duplicate Summary**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                  | Parameter                      | AOC 20-3 (0-0.5')<br>Sample Result | AOC 20-3 (0-0.5') DUP<br>Field Duplicate | RPD<br>(%) |
|------------------|--------------------------------|------------------------------------|--|------------|
| TPH (mg/kg-dry): | Diesel Range Organics (DRO)    | 87                                 | 150                                      | 53.1       |
|                  | Motor Oil Range Organics (MRO) | 78                                 | 120                                      | 42.2       |
|                  | Gasoline Range Organics (GRO)  | < 25                               | < 25                                     | NC         |
| VOCs (ug/kg-dry) | 1,1,1,2-Tetrachloroethane      | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1,1-Trichloroethane          | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1,2,2-Tetrachloroethane      | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1,2-Trichloroethane          | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1-Dichloroethane             | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1-Dichloroethene             | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,1-Dichloropropene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2,3-Trichlorobenzene         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2,3-Trichloropropane         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2,4-Trichlorobenzene         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2,4-Trimethylbenzene         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2-Dibromo-3-chloropropane    | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2-Dibromoethane (EDB)        | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2-Dichlorobenzene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,2-Dichloroethane (EDC)       | < 3.00                             | < 3.00                                   | NC         |
|                  | 1,2-Dichloropropane            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,3,5-Trimethylbenzene         | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,3-Dichlorobenzene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,3-Dichloropropane            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1,4-Dichlorobenzene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 1-Methylnaphthalene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 2,2-Dichloropropane            | < 1.00                             | < 1.00                                   | NC         |
|                  | 2-Butanone                     | < 1.00                             | < 1.00                                   | NC         |
|                  | 2-Chlorotoluene                | < 1.00                             | < 1.00                                   | NC         |
|                  | 2-Hexanone                     | < 1.00                             | < 1.00                                   | NC         |
|                  | 2-Methylnaphthalene            | < 1.00                             | < 1.00                                   | NC         |
|                  | 4-Chlorotoluene                | < 1.00                             | < 1.00                                   | NC         |
|                  | 4-Isopropyltoluene             | 4.77                               | < 1.00                                   | NC         |
|                  | 4-Methyl-2-pentanone           | < 1.00                             | < 1.00                                   | NC         |
|                  | Acetone                        | < 2.00                             | < 2.00                                   | NC         |
|                  | Benzene                        | < 1.00                             | < 1.00                                   | NC         |
|                  | Bromobenzene                   | < 1.00                             | < 1.00                                   | NC         |
|                  | Bromodichloromethane           | < 1.00                             | < 1.00                                   | NC         |
|                  | Bromoform                      | < 1.00                             | < 1.00                                   | NC         |
|                  | Bromomethane                   | < 1.00                             | < 1.00                                   | NC         |
|                  | Carbon disulfide               | 2.91                               | 4.61                                     | 45.2       |
|                  | Carbon tetrachloride           | < 1.00                             | < 1.00                                   | NC         |
|                  | Chlorobenzene                  | < 1.00                             | < 1.00                                   | NC         |
|                  | Chloroethane                   | < 1.00                             | < 1.00                                   | NC         |
|                  | Chloroform                     | < 1.00                             | < 1.00                                   | NC         |
|                  | Chloromethane                  | < 1.00                             | < 1.00                                   | NC         |
|                  | cis-1,2-DCE                    | < 1.00                             | < 1.00                                   | NC         |
|                  | cis-1,3-Dichloropropene        | < 1.00                             | < 1.00                                   | NC         |
|                  | Dibromochloromethane           | < 1.00                             | < 1.00                                   | NC         |
|                  | Dibromomethane                 | < 1.00                             | < 1.00                                   | NC         |
|                  | Dichlorodifluoromethane        | < 1.00                             | < 1.00                                   | NC         |
|                  | Ethylbenzene                   | < 1.00                             | < 1.00                                   | NC         |
|                  | Hexachlorobutadiene            | < 1.00                             | < 1.00                                   | NC         |
|                  | Isopropylbenzene               | < 1.00                             | < 1.00                                   | NC         |
|                  | Methyl tert-butyl ether (MTBE) | < 1.00                             | < 1.00                                   | NC         |
|                  | Methylene chloride             | 14.9                               | 13.1                                     | 12.8       |
|                  | Naphthalene                    | < 1.00                             | < 1.00                                   | NC         |
|                  | n-Butylbenzene                 | < 1.00                             | < 1.00                                   | NC         |
|                  | n-Propylbenzene                | < 1.00                             | < 1.00                                   | NC         |
|                  | sec-Butylbenzene               | < 1.00                             | < 1.00                                   | NC         |
|                  | Styrene                        | < 1.00                             | < 1.00                                   | NC         |
|                  | tert-Butylbenzene              | < 1.00                             | < 1.00                                   | NC         |
|                  | Tetrachloroethene (PCE)        | < 1.00                             | < 1.00                                   | NC         |
|                  | Toluene                        | 14.3                               | 15.8                                     | 9.9        |
|                  | trans-1,2-DCE                  | < 1.00                             | < 1.00                                   | NC         |
|                  | trans-1,3-Dichloropropene      | < 1.00                             | < 1.00                                   | NC         |
|                  | Trichloroethene (TCE)          | < 1.00                             | < 1.00                                   | NC         |
|                  | Trichlorofluoromethane         | < 1.00                             | < 1.00                                   | NC         |
|                  | Vinyl chloride                 | < 1.00                             | < 1.00                                   | NC         |
|                  | Xylenes, Total                 | < 1.00                             | < 1.00                                   | NC         |

**Table A-3**  
**Field Duplicate Summary**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                    | Parameter                   | AOC 20-3 (0-0.5') | AOC 20-3 (0-0.5') DUP | RPD (%) |
|--------------------|-----------------------------|-------------------|-----------------------|---------|
|                    |                             | Sample Result     | Field Duplicate       |         |
| SVOCs (mg/kg-dry): | 1,2,4-Trichlorobenzene      | < 0.40            | < 0.40                | NC      |
|                    | 1,2-Dichlorobenzene         | < 0.40            | < 0.40                | NC      |
|                    | 1,3-Dichlorobenzene         | < 0.40            | < 0.40                | NC      |
|                    | 1,4-Dichlorobenzene         | < 0.40            | < 0.40                | NC      |
|                    | 1-Methylnaphthalene         | < 0.40            | < 0.40                | NC      |
|                    | 2,4,5-Trichlorophenol       | < 0.40            | < 0.40                | NC      |
|                    | 2,4,6-Trichlorophenol       | < 0.40            | < 0.40                | NC      |
|                    | 2,4-Dichlorophenol          | < 0.80            | < 0.80                | NC      |
|                    | 2,4-Dimethylphenol          | < 0.60            | < 0.60                | NC      |
|                    | 2,4-Dinitrophenol           | < 0.80            | < 0.80                | NC      |
|                    | 2,4-Dinitrotoluene          | < 0.99            | < 1.0                 | NC      |
|                    | 2,6-Dinitrotoluene          | < 0.99            | < 1.0                 | NC      |
|                    | 2-Chloronaphthalene         | < 0.50            | < 0.50                | NC      |
|                    | 2-Chlorophenol              | < 0.40            | < 0.40                | NC      |
|                    | 2-Methylnaphthalene         | < 0.40            | < 0.40                | NC      |
|                    | 2-Methylphenol              | < 0.99            | < 1.0                 | NC      |
|                    | 2-Nitroaniline              | < 0.40            | < 0.40                | NC      |
|                    | 2-Nitrophenol               | < 0.40            | < 0.40                | NC      |
|                    | 3,3'-Dichlorobenzidine      | < 0.50            | < 0.50                | NC      |
|                    | 3+4-Methylphenol            | < 0.40            | < 0.40                | NC      |
|                    | 3-Nitroaniline              | < 0.40            | < 0.40                | NC      |
|                    | 4,6-Dinitro-2-methylphenol  | < 0.99            | < 1.0                 | NC      |
|                    | 4-Bromophenyl phenyl ether  | < 0.40            | < 0.40                | NC      |
|                    | 4-Chloro-3-methylphenol     | < 0.99            | < 1.0                 | NC      |
|                    | 4-Chloroaniline             | < 0.99            | < 1.0                 | NC      |
|                    | 4-Chlorophenyl phenyl ether | < 0.40            | < 0.40                | NC      |
|                    | 4-Nitroaniline              | < 0.80            | < 0.80                | NC      |
|                    | 4-Nitrophenol               | < 0.50            | < 0.50                | NC      |
|                    | Acenaphthene                | < 0.40            | < 0.40                | NC      |
|                    | Acenaphthylene              | < 0.40            | < 0.40                | NC      |
|                    | Aniline                     | < 0.40            | < 0.40                | NC      |
|                    | Anthracene                  | < 0.40            | < 0.40                | NC      |
|                    | Azobenzene                  | < 0.40            | < 0.40                | NC      |
|                    | Benz(a)anthracene           | < 0.40            | < 0.40                | NC      |
|                    | Benzo(a)pyrene              | < 0.40            | < 0.40                | NC      |
|                    | Benzo(b)fluoranthene        | < 0.40            | < 0.40                | NC      |
|                    | Benzo(g,h,i)perylene        | < 0.40            | < 0.40                | NC      |
|                    | Benzo(k)fluoranthene        | < 0.40            | < 0.40                | NC      |
|                    | Benzoic acid                | < 0.99            | < 1.0                 | NC      |
|                    | Benzyl alcohol              | < 0.40            | < 0.40                | NC      |
|                    | Bis(2-chloroethoxy)methane  | < 0.40            | < 0.40                | NC      |
|                    | Bis(2-chloroethyl)ether     | < 0.40            | < 0.40                | NC      |
|                    | Bis(2-chloroisopropyl)ether | < 0.40            | < 0.40                | NC      |
|                    | Bis(2-ethylhexyl)phthalate  | < 0.99            | < 1.0                 | NC      |
|                    | Butyl benzyl phthalate      | < 0.40            | < 0.40                | NC      |
|                    | Carbazole                   | < 0.40            | < 0.40                | NC      |
|                    | Chrysene                    | < 0.40            | < 0.40                | NC      |
|                    | Dibenz(a,h)anthracene       | < 0.40            | < 0.40                | NC      |
|                    | Dibenzofuran                | < 0.40            | < 0.40                | NC      |
|                    | Diethyl phthalate           | < 0.40            | < 0.40                | NC      |
|                    | Dimethyl phthalate          | < 0.40            | < 0.40                | NC      |
|                    | Di-n-butyl phthalate        | < 0.99            | < 1.0                 | NC      |
|                    | Di-n-octyl phthalate        | < 0.50            | < 0.50                | NC      |
|                    | Fluoranthene                | < 0.40            | < 0.40                | NC      |
|                    | Fluorene                    | < 0.40            | < 0.40                | NC      |
|                    | Hexachlorobenzene           | < 0.40            | < 0.40                | NC      |
|                    | Hexachlorobutadiene         | < 0.40            | < 0.40                | NC      |
|                    | Hexachlorocyclopentadiene   | < 0.40            | < 0.40                | NC      |
|                    | Hexachloroethane            | < 0.40            | < 0.40                | NC      |
|                    | Indeno(1,2,3-cd)pyrene      | < 0.40            | < 0.40                | NC      |
|                    | Isophorone                  | < 0.99            | < 1.0                 | NC      |
|                    | Naphthalene                 | < 0.40            | < 0.40                | NC      |
|                    | Nitrobenzene                | < 0.99            | < 1.0                 | NC      |
|                    | N-Nitrosodi-n-propylamine   | < 0.40            | < 0.40                | NC      |
|                    | N-Nitrosodiphenylamine      | < 0.40            | < 0.40                | NC      |
|                    | Pentachlorophenol           | < 0.80            | < 0.80                | NC      |
|                    | Phenanthrene                | < 0.40            | < 0.40                | NC      |
|                    | Phenol                      | < 0.40            | < 0.40                | NC      |
|                    | Pyrene                      | < 0.40            | < 0.40                | NC      |
|                    | Pyridine                    | < 0.99            | < 1.0                 | NC      |

**Table A-3**  
**Field Duplicate Summary**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                     | Parameter | AOC 20-3 (0-0.5') | AOC 20-3 (0-0.5') DUP | RPD (%) |
|---------------------|-----------|-------------------|-----------------------|---------|
|                     |           | Sample Result     | Field Duplicate       |         |
| Metals (mg/kg-dry): | Antimony  | < 13              | < 13                  | NC      |
|                     | Arsenic   | < 13              | < 13                  | NC      |
|                     | Barium    | 180               | 160                   | 11.7    |
|                     | Beryllium | 1.6               | 0.86                  | 60.1    |
|                     | Cadmium   | < 0.50            | < 0.50                | NC      |
|                     | Chromium  | 11                | 9.7                   | 12.5    |
|                     | Cobalt    | 8.9               | 8.0                   | 10.6    |
|                     | Cyanide   | < 0.30            | < 0.30                | NC      |
|                     | Lead      | 8.5               | 7.3                   | 15      |
|                     | Mercury   | < 0.033           | < 0.033               | NC      |
|                     | Nickel    | 9.0               | 8.7                   | 3.3     |
|                     | Selenium  | < 13              | < 13                  | NC      |
|                     | Silver    | < 1.3             | < 1.3                 | NC      |
|                     | Vanadium  | 32                | 24                    | 28.5    |
|                     | Zinc      | 42                | 42                    | 0.0     |

**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

\* = Field Duplicate RPD Outlier

**TABLE A-4**  
**Completeness Summaries**

**Table A-4**  
**Completeness Summary - Soil**  
**Group 6 Investigation Report**  
**Western Refining Southwest, Inc. - Bloomfield Refinery**

|                     | Parameter                      | Total Number of Results | Number of Contractual Compliance | Percent Contractual Compliance | Number of Usable Results | Percent Technical Compliance |
|---------------------|--------------------------------|-------------------------|----------------------------------|--------------------------------|--------------------------|------------------------------|
| TPH (mg/kg-dry):    | Diesel Range Organics (DRO)    | 45                      | 44 <sup>b</sup>                  | 97.8                           | 45                       | 100                          |
|                     | Motor Oil Range Organics (MRO) | 45                      | 45                               | 100.0                          | 45                       | 100                          |
|                     | Gasoline Range Organics (GRO)  | 45                      | 45                               | 100.0                          | 45                       | 100                          |
| VOCs (ug/kg-dry)    | 1-Methylnaphthalene            | 38                      | 38                               | 100.0                          | 38                       | 100                          |
|                     | 2-Methylnaphthalene            | 38                      | 38                               | 100.0                          | 38                       | 100                          |
|                     | 2-Butanone                     | 45                      | 44 <sup>a</sup>                  | 97.8                           | 45                       | 100                          |
|                     | Acetone                        | 45                      | 44 <sup>a</sup>                  | 97.8                           | 45                       | 100                          |
|                     | All remaining VOC analytes     | 45                      | 45                               | 100.0                          | 45                       | 100                          |
| SVOCs (mg/kg-dry):  | All SVOCs                      | 45                      | 45                               | 100                            | 45                       | 100                          |
| Metals (mg/kg-dry): | Antimony                       | 45                      | 26 <sup>c</sup>                  | 57.8                           | 45                       | 100                          |
|                     | Selenium                       | 45                      | 38 <sup>c</sup>                  | 84.4                           | 45                       | 100                          |
|                     | All remaining metals           | 45                      | 45                               | 100.0                          | 45                       | 100                          |

**Notes:**

Number of samples used in completeness calculations includes field duplicates but does not include equipment rinsate, field, or trip blanks.

Percent Contractual Compliance = (number of contract compliant results / Number of reported results)\*100

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

a = Qualified due to equipment blank detection.

b = Qualified due to high surrogate recoveries

c = Qualified due to low MS/MSD recovery



SUSANA MARTINEZ  
Governor

JOHN A. SANCHEZ  
Lieutenant Governor

NEW MEXICO  
ENVIRONMENT DEPARTMENT

*Hazardous Waste Bureau*

2905 Rodeo Park Drive East, Building 1  
Santa Fe, New Mexico 87505-6303  
Phone (505) 476-6000 Fax (505) 476-6030  
[www.nmenv.state.nm.us](http://www.nmenv.state.nm.us)



DAVE MARTIN  
Secretary

RAJ SOLOMON, P.E.  
Deputy Secretary

CERTIFIED MAIL - RETURN RECEIPT REQUESTED

March 1, 2011

Mr. Randy Schmaltz  
Environmental Manager  
Western Refining, Bloomfield Refinery  
P.O. Box 159  
Bloomfield, New Mexico 87413

**RE: APPROVAL  
INVESTIGATION WORK PLAN GROUP 6 (AOC NO. 19 SEEP  
NORTH OF MW-45, AOC NO. 20 SEEP NORTH OF MW-46, AND  
AOC NO. 21 SEEP NORTH OF MW-47)  
WESTERN REFINING SOUTHWEST, INC., BLOOMFIELD REFINERY  
EPA ID# NMD089416416  
HWB-WRB-10-002**

Dear Mr. Schmaltz:

The New Mexico Environment Department (NMED) has reviewed Western Refining Southwest, Inc., Bloomfield Refinery's (Western) *Investigation Work Plan Group 6 (AOC No. 19 Seep North of MW-45, AOC No. 20 Seep North of MW-46, and AOC No. 21 Seep North of MW-47)* (Work Plan), dated February 2011. NMED hereby approves this Work Plan with the following comments.

**Comment 1**

Figure 5 (Potentiometric Surface Map) and Figure 7 (Dissolved-Phase Groundwater Data) provide text boxes that report groundwater elevations and data results. In Figure 7, the groundwater data result boxes for MW-50 through MW-54 are easier to read than the other results included in Figure 7. In future documents, ensure that all text boxes are legible.



Mr. Schmaltz  
March 1, 2011  
Page 2 of 2

**Comment 2**

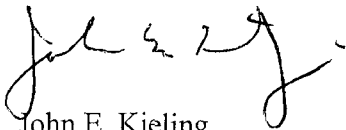
Figure 6 (Separate Phase Hydrocarbon Thickness Map) provides separate phase hydrocarbon (SPH) thickness measurements and contours for the whole refinery. The figure illustrates the light non-aqueous phase liquid (LNAPL) thickness with blue contours, but the corresponding LNAPL thickness measurements are reported in red font, which could be interpreted as representing a value other than LNAPL thickness. To avoid confusion in future documents, be consistent when reporting the SPH measurements and contours by keeping them the same color.

**Comment 3**

Appendix A (Photographs) includes photographs of all the seeps associated with the Group 6 AOCs. Photograph "AOC No. 19 – Seep No. 2" is clearly shown. However, it is difficult to view Seep No. 7 from the photograph "AOC 20 – Looking North Toward Seep No. 7 From Berm Below Catchment Basin No. 6." Future documents containing photographs of this nature must include arrows and labels on the photographs to clearly depict the locations of the seeps and catchment basins.

Western must adhere to all requirements in this Approval in future submittals. The Investigation Report must be submitted to NMED on or before November 15, 2011. If you have any questions regarding this letter, please contact Leona Tsinnajinnie of my staff at (505) 476-6057.

Sincerely,



John E. Kielling  
Program Manager  
Permits Management Program  
Hazardous Waste Bureau

cc: D. Cobrain, NMED HWB  
H. Monzeglio, NMED HWB  
C. Chavez, OCD  
A. Hains, Western

File: HWB-WRB-10-002 and Reading

RECEIVED

2009 DEC 30 PM 1 35

December 22, 2009

James Bearzi, Bureau Chief  
New Mexico Environmental Department  
Hazardous Waste Bureau  
2905 Rodeo Park Drive East, Building 1  
Santa Fe, New Mexico 87505-6303

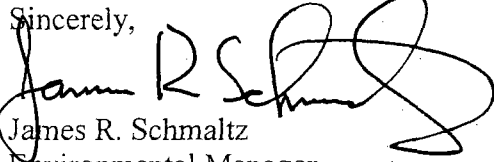
Re: Giant Refining Company, Bloomfield Refinery (currently know as Western Refinery Southwest, Inc. – Bloomfield Refinery) Order No. HWB 07-34 (CO) Investigation Work Plan Group No. 6

Dear Mr. Bearzi:

Western Refining Southwest Inc. - Bloomfield Refinery submits the referenced Investigation Work Plan pursuant to Section IV.B.4 of the July 2007 HWB Order. The Investigation Work Plan covers Group No. 6, which includes AOC No. 19 Seep North of MW-45, AOC No. 20 Seep North of MW-46, and AOC No. 21 Seep North of MW-47. The Investigation Work Plan was developed and formatted to meet the requirements of Section X.B of the July 2007 HWB Order.

If you have any questions or would like to discuss the Investigation Work Plan, please contact me at (505) 632-4171.

Sincerely,

  
James R. Schmaltz  
Environmental Manager  
Western Refining Southwest, Inc.  
Bloomfield Refinery

cc: Hope Monzeglio – NMED HWB  
Carl Chavez – NMOCD (w/attachment)  
Dave Cobrain – NMED HWB  
Laurie King – EPA Region 6 (w/attachment)  
Todd Doyle – Bloomfield Refinery  
Allen Hains – Western Refining El Paso

**RPS**

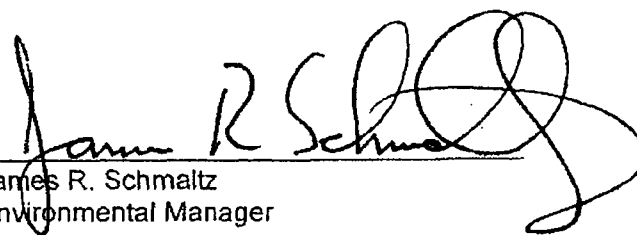
404 Camp Craft Rd., Austin, Texas 78746, USA  
T +1 512 347 7588 F +1 512 347 8243 W [www.rpsgroup.com](http://www.rpsgroup.com)

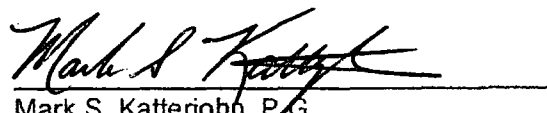
## **INVESTIGATION WORK PLAN Group 6**

(AOC No. 19 Seep North of MW-45,  
AOC No. 20 Seep North of MW-46,  
and AOC No. 21 Seep North of MW-47)

**Bloomfield Refinery  
Western Refining Southwest, Inc.  
#50 Rd 4990  
Bloomfield, New Mexico 87413**

**December 2009**

  
James R. Schmaltz  
Environmental Manager  
Western Refining Southwest, Inc.  
Bloomfield Refinery

  
Mark S. Katterjohn, P.G.  
Senior Consultant  
RPS  
404 Camp Craft Rd.  
Austin, Texas 78746

# Table of Contents

## List of Sections

|  |     |
|--|-----|
| Executive Summary .....  | iii |
| Section 1 Introduction.....  | 1   |
| Section 2 Background .....   | 2   |
| Section 3 Site Conditions .....                                    | 5   |
| 3.1 Surface Conditions .....                                       | 5   |
| 3.2 Subsurface Conditions.....                                     | 6   |
| Section 4 Scope of Services.....                                   | 7   |
| 4.1 Anticipated Activities.....                                    | 7   |
| 4.2 Background Information Research.....                           | 9   |
| 4.3 Collection and Management of Investigation Derived Waste ..... | 9   |
| 4.4 Surveys .....  | 9   |
| Section 5 Investigation Methods.....                               | 10  |
| 5.1 Soil Boring Installation .....                                 | 10  |
| 5.2 Soil Sampling .....  | 11  |
| 5.2.1 Soil Sample Field Screening and Logging .....                | 13  |
| 5.3 Sample Handling .....  | 14  |
| 5.4 Decontamination Procedures.....                                | 15  |
| 5.5 Field Equipment Calibration Procedures.....                    | 16  |
| 5.6 Documentation of Field Activities .....                        | 16  |
| 5.7 Chemical Analyses .....  | 17  |
| 5.8 Data Quality Objectives .....                                  | 18  |
| Section 6 Monitoring and Sampling Program.....                     | 20  |
| Section 7 Schedule .....   | 21  |
| Section 8 References .....   | 22  |

# **Table of Contents (*continued*)**

## **List of Tables**

|         |  |
|---------|--|
| Table 1 | Summary of 2007 Groundwater Seep Concentrations              |
| Table 2 | Summary of Facility-Wide Monitoring Plan Seep Concentrations |

## **List of Figures**

|           |  |
|-----------|--|
| Figure 1  | Site Location Map  |
| Figure 2  | AOC Group No. 6 Locations                                  |
| Figure 3  | Cross Section A-A' West to East                            |
| Figure 4  | Cross Section B-B' North to South                          |
| Figure 5  | Potentiometric Surface Map, August, 2008                   |
| Figure 6  | Separate Phase Hydrocarbon Thickness Map, August, 2008     |
| Figure 7  | Dissolved-Phase Groundwater Data, August and October, 2008 |
| Figure 8  | AOC Nos. 19, 20, and 21 Seep Location Map                  |
| Figure 9  | AOC No. 19 Sample Location Map                             |
| Figure 10 | AOC No. 20 Sample Location Map                             |
| Figure 11 | AOC No. 21 Sample Location Map                             |

## **List of Appendices**

|            |   |
|------------|---|
| Appendix A | Photographs                                       |
| Appendix B | Investigation Derived Waste (IDW) Management Plan |

## Executive Summary

The Bloomfield Refinery, which is located in the Four Corners Area of New Mexico, has been in operation since the late 1950s. Past inspections by State and federal environmental inspectors have identified locations where releases to the environment may have occurred. These locations are generally referred to as Solid Waste Management Units (SWMUs) or Areas of Concern (AOCs).

Pursuant to the terms and conditions of an Order issued on July 27, 2007 by the New Mexico Environment Department (NMED) to San Juan Refining Company and Giant Industries Arizona, Inc. for the Bloomfield Refinery, this Investigation Work Plan has been prepared for the AOCs designated as Group 6. A Class I modification to the facility's RCRA permit was approved on June 10, 2008 to reflect the change in ownership of the refinery to Western Refining Southwest, Inc. The operator is now Western Refining Southwest, Inc. – Bloomfield Refinery

The planned investigation activities include collection of soil samples, which will be analyzed for potential site-related constituents. The Investigation Work Plan includes specific sampling locations, sample collection procedures, and analytical methods. The scope of the proposed investigation is based, in part, on the results of previous site investigation activities.

Group 6 covers seeps located along the San Juan River northwest of the processing units, which are identified as AOC No. 19 Seep North of MW-45, AOC No. 20 Seep North of MW-46, and AOC No. 21 Seep North of MW-47 in the NMED Order. The Order requires that San Juan Refining Company and Giant Industries Arizona, Inc. ("Western") determine and evaluate the presence, nature, and extent of historical releases of contaminants at the aforementioned AOCs. This Investigation Work Plan has been developed to collect the necessary data to meet the requirements of the Order.

## Section 1

### Introduction

The Bloomfield Refinery is located immediately south of Bloomfield, New Mexico in San Juan County (Figure 1). The physical address is #50 Road 4990, Bloomfield, New Mexico 87413. The Bloomfield Refinery is located on approximately 263 acres. Bordering the facility is a combination of federal and private properties. Public property managed by the Bureau of Land Management lies to the south. The majority of undeveloped land in the vicinity of the facility is used extensively for oil and gas production and, in some instances, grazing. U.S. Highway 44 is located approximately one-half mile west of the facility. The topography of the main portion of the site is generally flat with steep bluffs to the north where the San Juan River intersects Tertiary terrace deposits.

The Bloomfield Refinery is a crude oil refinery currently owned by Western Refining Southwest, Inc., which is a wholly owned subsidiary of Western Refining Company, and it is operated by Western Refining Southwest, Inc. – Bloomfield Refinery. The Bloomfield Refinery has an approximate refining capacity of 18,000 barrels per day. Various process units are operated at the facility, including crude distillation, reforming, fluidized catalytic cracking, sulfur recovery, merox treater, catalytic polymerization, and diesel hydrotreating. Current and past operations have produced gasoline, diesel fuels, jet fuels, kerosene, propane, butane, naphtha, residual fuel, fuel oils, and LPG.

On July 27, 2007, the New Mexico Environment Department (NMED) issued an Order to San Juan Refining Company and Giant Industries Arizona, Inc. ("Western") requiring investigation and corrective action at the Bloomfield Refinery. This Investigation Work Plan has been prepared for the Areas of Concern (AOC) designated as Group 6 in the Order, which includes AOC No. 19 Seep North of MW-45, AOC No. 20 Seep North of MW-46, and AOC No. 21 Seep North of MW-47. The location of AOC Nos. 19, 20 and 21 are shown on Figure 2. Photographs of select locations around the seeps are included in Appendix A.

The purpose of the site investigation is to determine and evaluate the presence, nature, and extent of releases of contaminants in accordance with 20.4.1.500 New Mexico Administrative Code (NMAC) incorporating 40 Code of Federal Regulations (CFR) Section 264.101. The investigation activities will be conducted in accordance with Section IV of the Order.

## Section 2 Background

This section presents background information for the AOCs, including a review of historical waste management and product storage activities to identify the following:

- Type and characteristics of all waste and all potential contaminants handled in the subject AOCs;
- Known and possible sources of contamination;
- History of releases; and
- Known extent of contamination.

From 2001 to 2002, the Hammond Irrigation Ditch was lined with concrete. The ditch extends along the Facility's northern and western boundaries between the refinery and the San Juan River. The water present in the ditch acted as a hydraulic barrier prior to installation of the liner, confining the bulk of the contaminants to within the refinery boundaries. A recovery system with a French drain was installed to preserve the integrity of the concrete liner.

Between August and December 2004, petroleum hydrocarbons and water were observed to be seeping out of seven areas on the north side of the refinery near monitoring wells MW-45, MW-46, and MW-47. During inspections of these areas, hydrocarbon was observed to have stained the soil. Figure 2 shows the locations of these seeps.

Four seeps (Nos. 1, 2, 5, and 9) are located within two small drainage tributaries to the San Juan River at a location north of monitoring well MW-45. These seeps are referred to as AOC No. 19 Seep North of MW-45. Two seeps (Seep Nos. 6 and 7) were discovered along the bluff north of monitoring well MW-46 and are referred to as AOC No. 20 Seep North of MW-46. The last seep (No. 8) is located along the bluff above the San Juan River north of monitoring well MW-47 and is referred to as AOC No. 21 Seep North of MW-47.

Nine catchment basins were installed as temporary measures when the seeps were first identified in 2004. A liner was installed at each catchment basin with the exception of Seep No. 4. Initially, the water accumulating in the catchment basins was pumped to Tank 37 and then to the Refinery's wastewater treatment system. The basins were constructed by placing high density polyethylene (HDPE) on top of the exposed bedrock after the visibly impacted soils had



been removed and earthen berms formed. The design of the catchment basins also allows the accumulation of rain water in addition to groundwater discharge.

The North Boundary Barrier Wall (slurry wall) was designed and installed by April 2005 as a permanent control measure to prevent further potential discharges of groundwater to surface water and the generation of seeps along the bluff. Subsequent to the installation of the slurry wall and with the continued recovery of groundwater from the Hammond Ditch French drain, the discharge from the seeps was significantly reduced.

Only Seep Nos. 1, 6, 7, 8, and 9 currently have any actual discharge of groundwater. Seep No. 1 has the most flow at the rate of a trickle. Seep No. 6 has about two thirds the flow of Seep No. 1 with Seep Nos. 7, 8, and 9 measured by drips with no continuous flow. Seep Nos. 2 and 5 did have historic flows, but groundwater discharge has ceased subsequent to the installation of the slurry wall.

Water samples were collected from the seeps in 2007. The samples were collected on three occasions (October 9, 2007, November 8, 2007, and December 10, 2007) from the seeps, which had an active discharge of groundwater. The samples were collected from Seep Nos. 1, 6, 7, 8, and 9 shown on Figure 8 and were identified on the chain-of-custodies and analytical reports as Outfall #1, Outfall #6, Outfall #7, Outfall #8, and Outfall #9. The analytical results are summarized in Table 1. All of the analyses were non-detect for BTEX (benzene, ethylbenzene, toluene, and xylenes) constituents, with one exception. The October sample from Seep No. 1 (Outfall #1) indicated very low concentrations of benzene and xylenes at 0.0029 mg/l and 0.0039 mg/l, respectively. Both subsequent samples from this location were non-detect for all BTEX constituents. All of the detected results and non-detect quantitation limits are less than the potentially applicable groundwater and surface water quality standards shown in Table 1.

Water samples have been collected from the seeps on a semiannual basis starting in 2008. As part of the Facility-Wide Groundwater Monitoring Plan, Western checks all of the identified seep locations (Nos. 1 through 9) on a bi-weekly basis to determine if there is evidence of an active groundwater discharge. The presence of groundwater discharges is recorded and water samples are collected during the semiannual sampling event, when water is present, from all seeps that have indicated an active groundwater discharge within the past six months. The analytical results from these seep samples are summarized in Table 2. No semivolatile organic compounds were detected in any of the seep samples and, therefore, are not included in the

summary table. All analytical results are included in the Groundwater Remediation and Monitoring Report, which is submitted to the NMED in April of each year.

Results of the historical analyses of seep samples show that chloride and sulfate are detected at concentrations above the New Mexico Water Quality Control Commission Standards for Groundwater. All of the 2008 and 2009 analyses were non-detect for BTEX (benzene, ethylbenzene, toluene, and xylenes). Methyl tert-butyl ether (MTBE) was sporadically reported in some seep samples at concentrations ranging from less than 0.001 to 0.048 mg/L.

## Section 3

### Site Conditions

The surface and subsurface conditions that could affect the fate and transport of any contaminants are discussed below. This information is based on recent visual observations and historical subsurface investigations.

#### 3.1 Surface Conditions

Regionally, the surface topography slopes toward the floodplain of the San Juan River, which runs along the northern boundary of the refinery complex. To the south of the refinery, the drainage is to the northwest. North of the refinery, across the San Juan River, surface water flows in a southeasterly direction toward the San Juan River. The active portion of the refinery property, where the process units and storage tanks are located, is generally of low relief with an overall northwest gradient of approximately 0.02 ft/ft. The refinery sits on an alluvial floodplain terrace deposit and there is a steep bluff (approx. drop of 90 feet) at the northern boundary of the refinery where the San Juan River intersects the floodplain terrace, which marks the southern boundary of the floodplain.

There are two locally significant arroyos, one immediately east and another immediately west of the refinery. These arroyos collect most of the surface water flows in the area, thus significantly reducing surface water flows across the refinery. A minor drainage feature is located on the eastern portion of the refinery, where the former Landfill Pond (SWMU No. 9) was located and there are several steep arroyos along the northern refinery boundary that capture local surface water flows and minor groundwater discharges.

The refinery complex is bisected by County Rd #4990 (Sullivan Road), which runs east-west. The process units, storage tanks (crude oil and liquid products), and wastewater treatment systems are located north of the county road. The crude oil and product loading racks, LPG storage tanks and loading racks, maintenance buildings/90-day storage area, pipeline offices, transportation truck shop, and the Class I injection well are located south of the county road. There is very little vegetation throughout these areas with most surfaces composed of concrete, asphalt, or gravel. The area between the refinery and the San Juan River does have limited vegetation on steep slopes that do not support dense vegetation.

### 3.2 Subsurface Conditions

Numerous soil borings and monitoring wells have been completed across the refinery property during previous site investigations and installation of the slurry wall, which runs along the northern and western refinery boundary. Based on the available site-specific and regional subsurface information, the site is underlain by the Quaternary Jackson Lake terrace deposits, which unconformably overlie the Tertiary Nacimiento Formation. The Jackson Lake deposits consist of fine grained sand, silt and clay that grades to coarse sand, gravel and cobble size material closer to the contact with the Nacimiento Formation. The Jackson Lake Formation is over 40 feet thick near the southeast portion of the site and generally thins to the northwest toward the San Juan River. The Nacimiento Formation is primarily composed of fine grained materials (e.g., carbonaceous mudstone/claystone with interbedded sandstones) with a reported local thickness of approximately 570 feet (Groundwater Technology Inc., 1994).

Figures 3 and 4 present cross-sections of the shallow subsurface based on borings logs from on-site monitoring well completions. The uppermost aquifer is under water table conditions and occurs within the sand and gravel deposits of the Jackson Lake Formation. The Nacimiento Formation functions as an aquitard at the site that prevents contaminants from migrating to deeper aquifers. The potentiometric surface as measured in August 2008 is presented as Figure 5 and shows the groundwater flowing to the northwest, toward the San Juan River.

Previous investigations have identified and delineated impacts to groundwater from historical site operations. Figure 6 shows the distribution of separate phase hydrocarbon (SPH) in the subsurface based on the apparent thickness of SPH measured in monitoring wells. Dissolved-phase impacts are depicted on Figure 7.

## Section 4

### Scope of Services

#### 4.1 Anticipated Activities

Pursuant to Section IV of the Order, a scope of services has been developed to determine and evaluate the presence, nature, extent, fate, and transport of contaminants. To accomplish this objective, surface and shallow soil samples will be collected at AOC No. 19 Seep North of MW-45, AOC No. 20 Seep North of MW-46, and AOC No. 21 Seep North of MW-47. No groundwater samples are planned as the water from the seeps is already collected and analyzed on a semiannual basis as part of the Facility-Wide Groundwater Monitoring Plan. Therefore, the site investigation will be limited to the collection of soil samples. Where the HDPE liners are present within the catchment basins, surface soil samples will be collected as discussed below and in Section 5.2. Hand augured soil borings will be installed at select locations along the arroyos (drainage pathways) and samples will be collected from the following intervals:

- 0-6";
- 18-24";
- The sample from each soil boring with the greatest apparent degree of contamination, based on field observations and field screening; and
- Any additional intervals as determined based on field screening results.

Six soil borings are proposed for locations in the middle of the drainage pathways and at locations just downstream from the berms of the catchment basins. Thirty four surface soil samples (0-6") will be collected at the seep faces, on top of the HDPE liner, at select locations along the bottom of the drainage pathways, and along the margins of the drainage pathways, which occur on the banks of the drainage pathways and at the edge of the catchment basin liners.

Soil samples will be collected at AOC No. 19 from the following locations:

- Three surface soil samples will be collected near Seep No. 1, with one of the samples collected from the seep face and the other two samples collected along the margins of the drainage pathway;

- Near Seep No.2, a total of three surface soil samples will be collected with one collected from the seep face and the other two collected from along the margins of the drainage pathway;
- Three surface soil samples will be collected near Seep No. 5 with one collected from the seep face and two more collected along the margins of the drainage pathway;
- Four surface soil samples will be collected near Seep No. 9 with one collected from the bottom of the drainage pathway where the seep originated, one from the center of the lined catchment basin, and two from along the margins of the drainage pathway;
- One surface soil sample will be collected at Seep No. 4 at the center of the natural basin where the seep daylighted;
- Two surface soil samples will be collected from along the margins of the drainage pathway between Seep No. 5 and Seep No. 9;
- Four surface soil samples will be collected along the margins of the drainage pathway between Seep No. 3 and Seep No. 4; and
- A minimum of six soil samples will be collected from the three soil borings, which are located near the center of the two major drainage pathways that occur at AOC No. 19.

At AOC No. 20, soil samples will be collected at the following locations:

- Four surface samples will be collected at Seep No. 6 with one sample collected from the seep face, one soil sample collected from on top of the catchment liner, and two samples collected from along the margins of the drainage pathway;
- Three surface soil samples will be collected at Seep No. 7, with one sample collected from the top of the catchment liner and two samples collected from along the margins of the drainage pathway;
- Two surface soil samples will be collected from along the center line of the drainage pathway that extends from Seep No. 6 toward the San Juan River;
- Two surface soil samples will be collected from the margins of the drainage pathway that extends from Seep No. 6 toward the San Juan River; and
- A minimum of two soil samples will be collected from one soil boring at the center line of the drainage pathway that extends below Seep No. 6.

Soil samples will be collected from the following locations at AOC No. 21:

- Three surface soil samples will be collected at Seep No. 8, with one collected from the top of the catchment liner and two collected from along the margins of the drainage pathway; and
- A minimum of four soil samples will be collected from two soil borings, with one positioned at the discharge point of Seep No. 8 and the other down the drainage pathway just below the catchment berm near Seep No. 8.

The soil samples will be analyzed by the following methods:

- SW-846 Method 8260 (volatile organic compounds);
- SW-846 Method 8270 (semi-volatile organic compounds);
- SW-846 Method 8015B gasoline range (C5-C10), diesel range (>C10-C28), and motor oil range (>C28-C36) organics;
- SW-846 Methods 6010/6020 and 7470/7471 (metals); and
- SW-846 Methods 335.4/335.2 (cyanide).

## **4.2 Background Information Research**

Documents containing the results of previous investigations and subsequent routine groundwater monitoring data from monitoring wells and the seeps were reviewed to facilitate development of this work plan. The previously collected data provide detailed information on the overall subsurface conditions, including hydrogeology and contaminant distribution within groundwater on a site-wide basis. The data collected under this scope of services will supplement the existing soil and groundwater information and provide specific information regarding contaminant occurrence and distribution within soils near the seeps.

## **4.3 Collection and Management of Investigation Derived Waste**

Excess sample material and decontamination fluids, and all other investigation derived waste (IDW) associated with site investigation activities will be contained and characterized using methods based on the boring location, boring depth, and type of contaminants suspected or encountered. All decontamination water will be disposed in the refinery wastewater treatment system upstream of the API Separator. An IDW management plan is included as Appendix B.

## **4.4 Surveys**

The horizontal coordinates and elevation of each sampling location and the locations of all other pertinent structures will be determined by a registered New Mexico professional land surveyor in accordance with the State Plane Coordinate System (NMSA 1978 47-1-49-56 (Repl. Pamp. 1993)). The surveys will be conducted in accordance with Sections 500.1 through 500.12 of the Regulations and Rules of the Board of Registration for Professional Engineers and Surveyors Minimum Standards for Surveying in New Mexico. Horizontal positions will be measured to the nearest 0.1-ft and vertical elevations will be measured to the nearest 0.01-ft.

## Section 5

# Investigation Methods

The purpose of the site investigation is to determine and evaluate the presence, nature, and extent of releases of contaminants. Guidance on selecting and developing sampling plans as provided in *Guidance for Choosing a Sampling Design for Environmental Data Collection* (EPA, 2000) was utilized to select the appropriate sampling strategy.

### 5.1 Soil Boring Installation

Shallow soil borings will be completed using a hand auger and/or shovel at appropriate locations within and into the banks of the drainage pathways where the HDPE liner of the catchment basins is not present. This sampling method is ideally suited due to the inaccessibility of a drill rig and the limited anticipated depth to bedrock in the vicinity of the seeps. The sampling equipment will be properly decontaminated before beginning each boring.

The NMED will be notified as early as practicable if conditions arise or are encountered that do not allow the advancement of borings at planned sampling locations. Appropriate actions (e.g., relocation of borings to a less threatening location) will be taken to minimize any negative impacts from investigative borings. The depth of soil borings in the vicinity of the seeps may be limited due to collapse of the overlying soil. The soil borings within the drainage pathways will be extended to the top of bedrock (Nacimiento Formation) or to refusal, whichever occurs first, at an anticipated completion depth of less than 3 feet. Soil samples will be collected continuously and logged by a qualified geologist or engineer.

The sampling will be accomplished under the direction of a qualified engineer or geologist who will maintain a detailed log of the materials and conditions encountered in each boring. Both sample information and visual observations of the cuttings and core samples will be recorded on the boring log. Known site features and/or site survey grid markers will be used as references to locate each boring prior to surveying the location as described in Section 4.4. The boring locations will be measured to the nearest foot, and locations will be recorded on a scaled site map upon completion of each boring.



## 5.2 Soil Sampling

Since there is the potential for constituents to have been released to soils at known locations at AOC Nos. 19, 20 and 21, a judgmental sampling design is appropriate. Soil samples will be collected from the locations where groundwater reaches the land surface (i.e., seep face) and along the bottom and sides of the drainage pathways where water flowed and accumulated. The individual sample locations have been selected in an attempt to determine the highest concentrations of constituents in soil and to determine the horizontal extent of the impacts. The surface soil samples (0-6") collected from the seep faces and soil that has accumulated on top of the catchment basin liners where water accumulated should represent the higher concentration areas. Samples collected from soil borings located at seep faces and along the center of the drainage pathways will also represent higher concentration areas. The drainage pathways are generally narrow and deeply incised into the alluvial terrace deposits overlying the bedrock (Nacimiento Formation). Surface soil samples (0-6") collected from the sides of the steep slopes along the drainage pathways and at the edge of catchment liners will be used to define the horizontal extent of any impacts identified in the seep faces and bottom of the drainage pathways.

The six soil borings will be completed using a hand auger to the top of bedrock or to refusal, whichever occurs first. The thirty four surface soil sample locations will terminate at the top of the HDPE-liner, where present, or at a depth of 6", unless Western elects to extend the sampling deeper based on field screening results. A decontaminated hand auger will be used to obtain disturbed soils during the installation of each boring. Surface samples may be collected using decontaminated, hand-held stainless steel sampling device, hand auger, or a pre-cleaned disposable sampling device. A portion of the sample will be placed in pre-cleaned, laboratory-prepared sample containers for laboratory chemical analysis. The use of an Encore® Sampler or other similar device will be used during collection of soil samples for VOC analysis. The remaining portions of the sample will be used for logging and field screening as discussed in Section 5.2.1. Sample handling and chain-of-custody procedures will be in accordance with the procedures presented below in Section 5.3.

Discrete soil samples will be collected at the soil borings for laboratory analyses at the following intervals:

- 0-6" (all soil sample locations);
- 18-24" (only from the six deeper hand-augured soil boring locations);

- The sample from each of the six hand-augured soil borings with the greatest apparent degree of contamination, based on field observations and field screening; and
- Any additional intervals as determined based on field screening results.

Proposed sampling locations for AOC Nos. 19, 20 and 21 are shown on Figures 9, 10 and 11, respectively and are described in detail below.

#### AOC No. 19

Soil samples will be collected at AOC No. 19 from the following locations:

- Three surface soil samples will be collected near Seep No. 1, with one of the samples collected from the seep face and the other two collected along the margins of the drainage pathway;
- Near Seep No.2, a total of three surface soil samples will be collected with one from the seep face and the other two from along the margins of the drainage pathway;
- Three surface soil samples will be collected near Seep No. 5 with one collected from the seep face and two more collected along the margins of the drainage pathway;
- Four surface soil samples will be collected near Seep No. 9 with one collected from the bottom of the drainage pathway where the seep originated, one from the center of the lined catchment basin, and two from along the margins of the drainage pathway;
- One surface soil sample will be collected at Seep No. 4 at the center of the natural basin where the seep daylighted;
- Two surface soil samples will be collected along the margins of the drainage pathway between Seep No. 5 and Seep No. 9;
- Four surface soil samples will be collected along the margins of the drainage pathway between Seep No. 3 and Seep No. 4; and
- A minimum of six soil samples will be collected from the three soil borings, which are located near the center of the two major drainage pathways that occur at AOC No. 19.

#### AOC No. 20

At AOC No. 20, soil samples will be collected at the following locations:

- Four surface samples will be collected at Seep No. 6 with one sample collected from the seep face, one soil sample collected from on top of the catchment liner, and two samples collected from along the margins of the drainage pathway;

- Three surface soil samples will be collected at Seep No. 7, with one sample collected from the top of the catchment liner and two samples collected from along the margins of the drainage pathway;
- Two surface soil samples will be collected from along the center line of the drainage pathway that extends from Seep No. 6 toward the San Juan River;
- Two surface soil samples will be collected from the margins of the drainage pathway that extends from Seep No. 6 toward the San Juan River; and
- A minimum of two soil samples will be collected from one soil boring at the center line of the drainage pathway that extends below Seep No. 6.

#### AOC No. 21

Soil samples will be collected from the following locations at AOC No. 21:

- Three surface soil samples will be collected at Seep No. 8, with one collected from the top of the catchment liner and two collected from along the margins of the drainage pathway; and
- A minimum of four soil samples will be collected from two soil borings, with one positioned at the discharge point of Seep No. 8 and another down the drainage pathway just below the catchment berm near Seep No. 8.

Quality Assurance/Quality Control (QA/QC) samples will be collected to monitor the validity of the soil sample collection procedures as follows:

- Field duplicates will be collected at a rate of 10 percent;
- Equipment blanks will be collected from all sampling apparatus at a frequency of 10 percent or one per day if disposable sampling equipment is used; and
- Field blanks will be collected at a frequency of one per day.

#### **5.2.1 Soil Sample Field Screening and Logging**

Samples obtained from the borings will be screened in the field on 1.0 foot intervals for evidence of contaminants. Field screening results will be recorded on the exploratory boring logs and will be used to aid in the selection of soil samples for laboratory analysis. The primary screening methods include: (1) visual examination, (2) olfactory examination, and (3) headspace vapor screening for volatile organic compounds.

Visual screening includes examination of soil samples for evidence of staining caused by petroleum-related compounds. Headspace vapor screening targets volatile organic compounds

and involves placing a soil sample in a plastic sample bag or a foil sealed container allowing space for ambient air. The container will be sealed and then shaken gently to expose the soil to the air trapped in the container. The sealed container will be allowed to rest for a minimum of 5 minutes while vapors equilibrate. Vapors present within the sample bag's headspace will then be measured by inserting the probe of the instrument in a small opening in the bag or through the foil. The maximum value and the ambient air temperature will be recorded on the field boring log for each sample.

The monitoring instruments will be calibrated each day to the manufacturer's standard for instrument operation. A photo-ionization detector (PID) equipped with a 10.6 or higher electron volt (eV) lamp or a combustible gas indicator will be used for VOC field screening. Field screening results may be site- and boring-specific and the results may vary with instrument type, the media screened, weather conditions, moisture content, soil type, and type of contaminant. Conditions capable of influencing the results of field screening will be recorded on the field logs.

The physical characteristics of the samples (such as mineralogy, ASTM soil classification, moisture content, texture, color, presence of stains or odors, and/or field screening results), depth where each sample was obtained, method of sample collection, and other observations will be recorded in the field log by a qualified geologist or engineer. Detailed logs of each soil boring will be completed in the field by a qualified engineer or geologist. Additional information, such as the presence of water-bearing zones and any unusual or noticeable conditions encountered during drilling, will be recorded on the logs.

### **5.3 Sample Handling**

At a minimum, the following procedures will be used at all times when collecting samples during investigation, corrective action, and monitoring activities:

1. Neoprene, nitrile, or other protective gloves will be worn when collecting samples. New disposable gloves will be used to collect each sample;
2. All samples collected of each medium for chemical analysis will be transferred into clean sample containers supplied by the project analytical laboratory with the exception of soil, rock, and sediment samples obtained in Encore® samplers. Sample container volumes and preservation methods will be in accordance with the most recent standard EPA and industry accepted practices for use by accredited analytical laboratories. Sufficient sample volume will be obtained for the laboratory to complete the method-specific QC analyses on a laboratory-batch basis; and

3. Sample labels and documentation will be completed for each sample following procedures discussed below. Immediately after the samples are collected, they will be stored in a cooler with ice or other appropriate storage method until they are delivered to the analytical laboratory. Standard chain-of-custody procedures, as described below, will be followed for all samples collected. All samples will be submitted to the laboratory soon enough to allow the laboratory to conduct the analyses within the method holding times. At a minimum, all samples will be submitted to the laboratory within 48 hours after their collection.

Chain-of-custody and shipment procedures will include the following:

1. Chain-of-custody forms will be completed at the end of each sampling day, prior to the transfer of samples off site.
2. Individual sample containers will be packed to prevent breakage and transported in a sealed cooler with ice or other suitable coolant or other EPA or industry-wide accepted method. The drainage hole at the bottom of the cooler will be sealed and secured in case of sample container leakage.
3. Each cooler or other container will be delivered directly to the analytical laboratory.
4. Glass bottles will be separated in the shipping container by cushioning material to prevent breakage.
5. Plastic containers will be protected from possible puncture during shipping using cushioning material.
6. The chain-of-custody form and sample request form will be shipped inside the sealed storage container to be delivered to the laboratory.
7. Chain-of-custody seals will be used to seal the sample-shipping container in conformance with EPA protocol.
8. Signed and dated chain-of-custody seals will be applied to each cooler prior to transport of samples from the site.
9. Upon receipt of the samples at the laboratory, the custody seals will be broken, the chain-of-custody form will be signed as received by the laboratory, and the conditions of the samples will be recorded on the form. The original chain-of-custody form will remain with the laboratory and copies will be returned to the relinquishing party.
10. Copies of all chain-of-custody forms generated as part of sampling activities will be maintained on-site.

#### **5.4 Decontamination Procedures**

The objective of the decontamination procedures is to minimize the potential for cross-contamination. A designated decontamination area will be established for decontamination of sampling equipment.

Sampling or measurement equipment, including but not limited to, stainless steel sampling tools will be decontaminated in accordance with the following procedures or other methods approved by the Department before each sampling attempt or measurement:

1. Brush equipment with a wire or other suitable brush, if necessary or practicable, to remove large particulate matter;
2. Rinse with potable tap water;
3. Wash with nonphosphate detergent or other detergent approved by the Department (examples include Fantastik™, Liqui-Nox®);
4. Rinse with potable tap water; and
5. Double rinse with deionized water.

All decontamination solutions will be collected and stored temporarily as described in Section 4.3. Decontamination procedures and the cleaning agents used will be documented in the daily field log.

## **5.5 Field Equipment Calibration Procedures**

Field equipment requiring calibration will be calibrated to known standards, in accordance with the manufacturers' recommended schedules and procedures. At a minimum, calibration checks will be conducted daily, or at other intervals approved by the Department, and the instruments will be recalibrated, if necessary. Calibration measurements will be recorded in the daily field logs. If field equipment becomes inoperable, its use will be discontinued until the necessary repairs are made. In the interim, a properly calibrated replacement instrument will be used.

## **5.6 Documentation of Field Activities**

Daily field activities, including observations and field procedures, will be recorded in a field log book. The original field forms will be maintained at the facility. Copies of the completed forms will be maintained in a bound and sequentially numbered field file for reference during field activities. Indelible ink will be used to record all field activities. Photographic documentation of field activities will be performed, as appropriate. The daily record of field activities will include the following:

1. Site or unit designation;
2. Date;
3. Time of arrival and departure;
4. Field investigation team members including subcontractors and visitors;
5. Weather conditions;

6. Daily activities and times conducted;
7. Observations;
8. Record of samples collected with sample designations and locations specified;
9. Photographic log, as appropriate;
10. Field monitoring data, including health and safety monitoring;
11. Equipment used and calibration records, if appropriate;
12. List of additional data sheets and maps completed;
13. An inventory of the waste generated and the method of storage or disposal; and
14. Signature of personnel completing the field record.

## **5.7 Chemical Analyses**

All samples collected for laboratory analysis will be submitted to an accredited laboratory. The laboratory will use the most recent standard EPA and industry-accepted analytical methods for target analytes as the testing methods for each medium sampled. Chemical analyses will be performed in accordance with the most recent EPA standard analytical methodologies and extraction methods.

The soil samples will be analyzed by the following methods:

- SW-846 Method 8260 volatile organic compounds;
- SW-846 Method 8270 semi-volatile organic compounds; and
- SW-846 Method 8015B gasoline range (C5-C10), diesel range (>C10-C28), and motor oil range (>C28-C36) organics.

Soil samples will also be analyzed for the following metals and cyanide using the indicated analytical methods.

| 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       |
| 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       |

## 5.8 Data Quality Objectives

The Data Quality Objectives (DQOs) were developed to ensure that newly collected data are of sufficient quality and quantity to address the projects goals, including Quality Assurance/Quality Control (QA/QC) issues (EPA, 2006). The project goals are established in the Order and are to determine and evaluate the presence, nature, and extent of releases of contaminants at specified AOCs. The type of data required to meet the project goals includes chemical analyses of soil, sediment and groundwater to determine if there has been a release of contaminants at the individual AOCs.

The quantity of data is AOC specific and based on the historical operations at individual locations. The quality of data required is consistent across locations and is specified in Section VIII.D.7.c of the Order. In general, method detection limits should be 20% or less of the applicable background levels, cleanup standards and screening levels.

Additional DQOs include precision, accuracy, representativeness, completeness, and comparability. Precision is a measurement of the reproducibility of measurements under a given set of circumstances and is commonly stated in terms of standard deviation or coefficient



of variation (EPA, 1987). Precision is also specific to sampling activities and analytical performance. Sampling precision will be evaluated through the analyses of duplicate field samples and laboratory replicates will be utilized to assess laboratory precision.

Accuracy is a measurement in the bias of a measurement system and may include many sources of potential error, including the sampling process, field contamination, preservation, handling, sample matrix, sample preparation, and analysis techniques (EPA, 1987). An evaluation of the accuracy will be performed by reviewing the results of field/trip blanks, matrix spikes, and laboratory QC samples.

Representativeness is an expression of the degree to which the data accurately and precisely represent the true environmental conditions. Sample locations and the number of samples have been selected to ensure the data is representative of actual environmental conditions. Based on AOC specific conditions, this may include either biased (i.e., judgmental) locations/depths or unbiased (systematic grid samples) locations, as discussed in Section 5.2 for soils. In addition, sample collection techniques (e.g., field monitoring and decontamination of sampling equipment) will be utilized to help ensure representative results.

Completeness is defined as the percentage of measurements taken that are actually valid measurements, considering field QA and laboratory QC problems. EPA Contract Laboratory Program (CLP) data has been found to be 80-85% complete on a nationwide basis and this has been extrapolated to indicate that Level III, IV, and V analytical techniques will generate data that are approximately 80% complete (EPA, 1987). As an overall project goal, the completeness goal is 85%; however, some samples may be critical based on location or field screening results and thus a sample-by-sample evaluation will be performed to determine if the completeness goals have been obtained.

Comparability is a qualitative parameter, which expresses the confidence with which one data set can be compared to another. Industry standard sample collection techniques and routine EPA analytical methods will be utilized to help ensure data are comparable to historical and future data. Analytical results will be reported in appropriate units for comparison to historical data and cleanup levels.

## **Section 6**

### **Monitoring and Sampling Program**

Water samples from the seeps are collected on a semiannual basis as part of the NMED-approved Facility-Wide Groundwater Monitoring Plan. In accordance with Section 5.3 of the Facility-Wide Groundwater Monitoring Plan, water samples are collected from the seeps, if water is present. All seep locations are visually inspected bi-weekly to monitor active groundwater discharge along the bluff (NMED, 2008). The analytical results are included in the Groundwater Remediation and Monitoring Report submitted in April of each year. Groundwater is removed from any seep where analytical results exceed any of the standards set by the Water Quality Control Commission (WQCC), the EPA Maximum Contaminant Level (MCL), or the EPA Region VI Human Health Medium Specific Screening Levels (Tap Water) in the absence of a WQCC standard or MCL (NMED, 2008).

## **Section 7 Schedule**

This investigation Work Plan will be implemented within 90 days of NMED approval. The estimated timeframes for each of the planned activities is as shown below:

- Field work (inclusive of all soil sampling) -- four weeks;
- Laboratory analyses for initial sampling event – eight weeks;
- Data reduction and validation (soils) – three weeks; and
- Data gap analysis – three weeks.

Completion of the data gap analysis will complete all activities conducted under this investigation Work Plan. Western will then prepare an Investigation Report pursuant to Section X.C of the Order. The Investigation Report will be submitted to the NMED within 120 calendar days of completion of the data gap analysis.

## Section 8

### References

- EPA, 1987, Data Quality Objectives for Remedial Response Activities; United States Environmental Protection Agency, Office of Emergency and Remedial Response and Office of Waste Programs Enforcement, OSWER Directive 9355.0-7B, 85p
- EPA, 2000, Guidance on Choosing a Sampling Design for Environmental Data Collection, EPA/240/R-02/005, EPA QA/G-5S, 168 p.
- EPA, 2006, Guidance on Systematic Planning Using the Data Quality Objectives Process, United States Environmental Protection Agency, Office of Environmental Information; EPA/240/B-06/001, p. 111.
- Groundwater Technology Inc., 1994, RCRA Facility Investigation/Corrective Measures Study Report Bloomfield Refining Company #50 County Road 4990 Bloomfield, New Mexico, p.51.
- NMED, 2007, State of New Mexico Environment Department v. San Juan Refining Company and Giant Industries, Inc.; Order July 27, 2007, p. 133
- NMED, 2008. New Mexico Environmental Department (NMED) Hazardous Waste Bureau, Approval with Direction Facility-Wide Groundwater Monitoring Plan dated March 25, 2008.

# Tables

---

**Table 1**  
**Summary of 2007 Groundwater Seep Concentrations**  
**Group 6 Investigation Work Plan**  
**Western Refining Southwest - Bloomfield Refinery**

|                             |               | Parameters        |                   |                        |                   |
|-----------------------------|---------------|-------------------|-------------------|------------------------|-------------------|
|                             |               | Benzene<br>(mg/L) | Toluene<br>(mg/L) | Ethylbenzene<br>(mg/L) | Xylenes<br>(mg/L) |
| WQCC 20NMAC 6.2.3103(mg/L): |               | 0.01              | 0.75              | 0.75                   | 0.62              |
| WQCC 20NMAC 6.4(mg/L):      |               | 0.022             | 6.8               | 3.1                    | NE                |
| Sample ID:                  | Date Sampled: |                   |                   |                        |                   |
| Outfall #1<br>(Seep No. 1)  | 10/9/2007     | <0.001            | 0.0029            | <0.001                 | 0.0039            |
|                             | 11/8/2007     | <0.001            | <0.001            | <0.001                 | <0.002            |
|                             | 12/10/2007    | <0.001            | <0.001            | <0.001                 | <0.002            |
| Outfall #6<br>(Seep No. 6)  | 10/9/2007     | <0.001            | <0.001            | <0.001                 | <0.002            |
|                             | 11/8/2007     | <0.001            | <0.001            | <0.001                 | <0.002            |
|                             | 12/10/2007    | <0.001            | <0.001            | <0.001                 | <0.002            |
| Outfall #7<br>(Seep No. 7)  | 10/9/2007     | <0.001            | <0.001            | <0.001                 | <0.002            |
|                             | 11/8/2007     | <0.001            | <0.001            | <0.001                 | <0.002            |
|                             | 12/10/2007    | <0.001            | <0.001            | <0.001                 | <0.002            |
| Outfall #8<br>(Seep No. 8)  | 10/9/2007     | <0.001            | <0.001            | <0.001                 | <0.002            |
|                             | 11/7/2007     | <0.001            | <0.001            | <0.001                 | <0.002            |
|                             | 12/10/2007    | <0.001            | <0.001            | <0.001                 | <0.002            |
| Outfall #9<br>(Seep No. 9)  | 10/9/2007     | <0.001            | <0.001            | <0.001                 | <0.002            |
|                             | 11/8/2007     | <0.001            | <0.001            | <0.001                 | <0.002            |
|                             | 12/10/2007    | <0.001            | <0.001            | <0.001                 | <0.002            |

**Notes:**

mg/L = milligram per liter

NE = not established

WQCC 20NMAC 6.2.3103 = New Mexico Standard for Groundwater of 10,000 ug/L or less

WQCC 20NMAC 6.4 = New Mexico Standards for Interstate and Intrastate Surface Water

**Table 2**  
**Summary of Facility-Wide Monitoring Plan Seep Concentrations**  
**Group 6 Investigation Work Plan**  
**Western Refining Southwest - Bloomfield Refinery**

| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
|---|--|--|--|--|--|--|--|--|--|--|--|--|--|--|
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| WQCC 20 NMAC 6.2.3103   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
| 40CFR141.61 (Benzene and Ethylbenzene) EPA Region VI Screening Level (MTBE) |  |  |  |  |  |  |  |  |  |  |  |  |  |  |
|   |  |  |  |  |  |  |  |  |  |  |  |  |  |  |

NS<sup>1</sup> = Seep location is Dry or Not Enough Water to Sample- No Sample

NS<sup>2</sup> = Not Sampled - Sample was taken before implementation of Facility-Wide Monitoring Plan

NA = Constituent analysis not requested

WQCC 20NMAC 6.2.3103 = New Mexico Water Quality Control Commission Standard for Groundwater of 10,000 ug/L or less

40CFR141.61 = Safe Drinking Water Act Maximum Contaminant Level

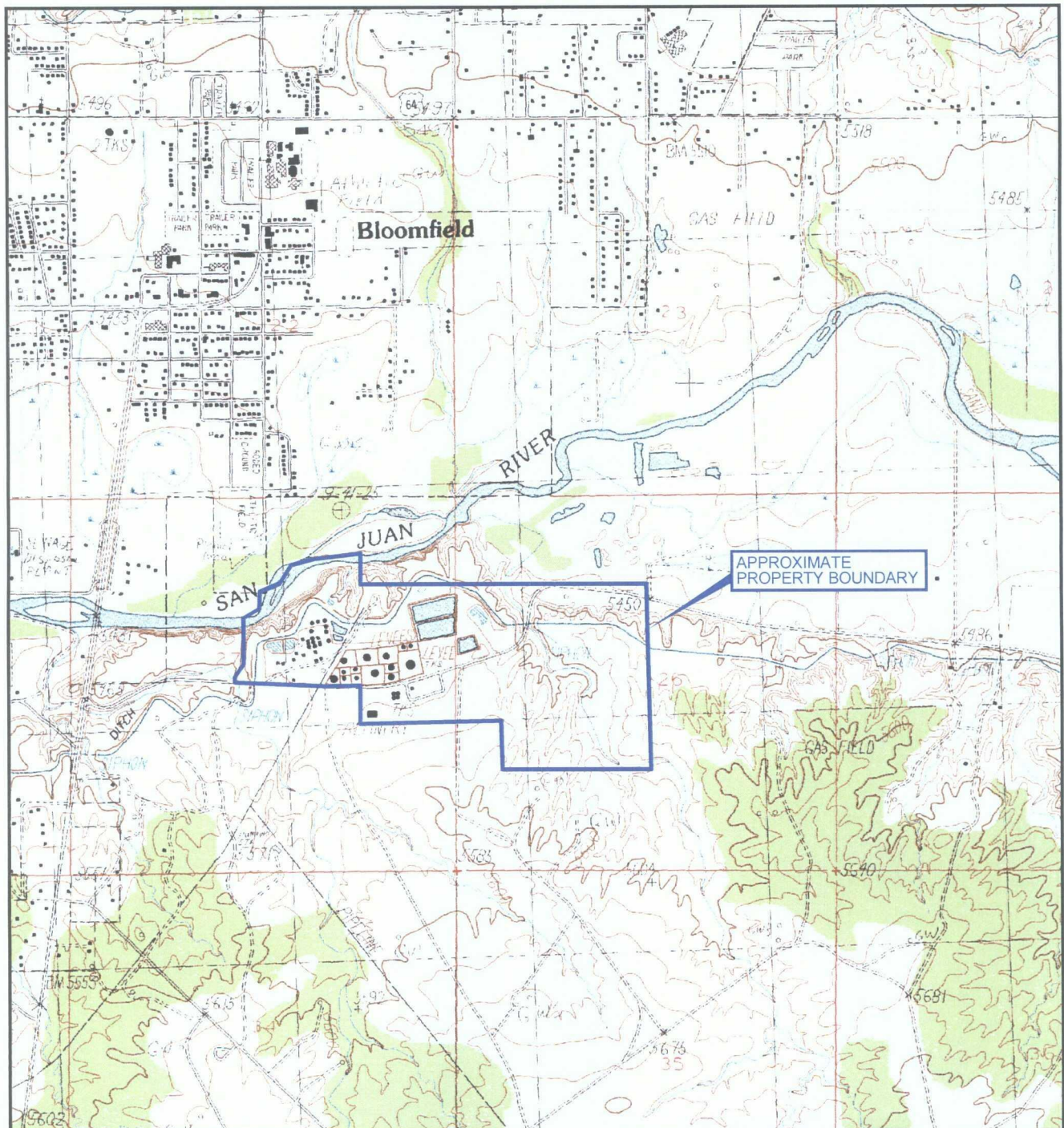
EPA Region VI Human Health Medium-Specific Screening Level - Tap Water

No active discharge at Seeps 2, 4 and 5 has been present to sample since implementation of the Facility-Wide Monitoring Plan

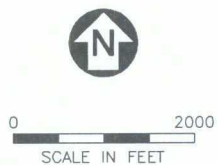
## Figures

---





Map Source: USGS 7.5 Min. Quad Sheet BLOOMFIELD, NM., 1985.



WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 6/19/08 FILE: WestRef-A25

FIGURE 1  
SITE LOCATION MAP  
BLOOMFIELD REFINERY



404 Camp Craft Road  
Austin, Texas 78746



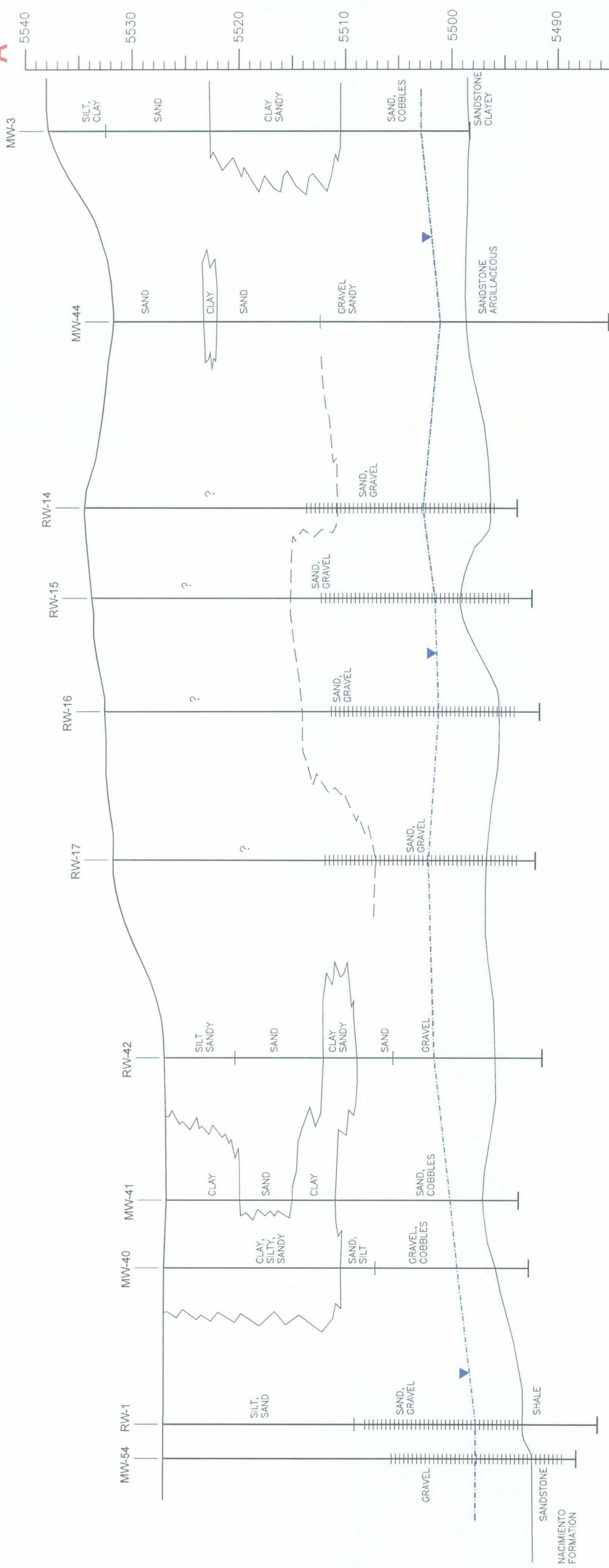




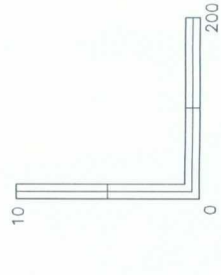
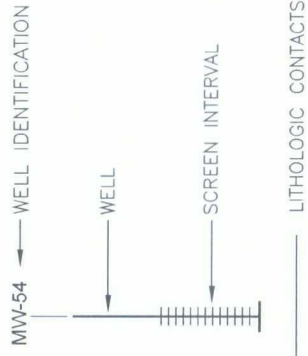
WEST  
A



EAST  
A'



EXPLANATION



**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 04/29/09 FILE: WestRef-B28

FIGURE 3  
CROSS SECTION A-A'  
WEST TO EAST  
BLOOMFIELD REFINERY

RPS

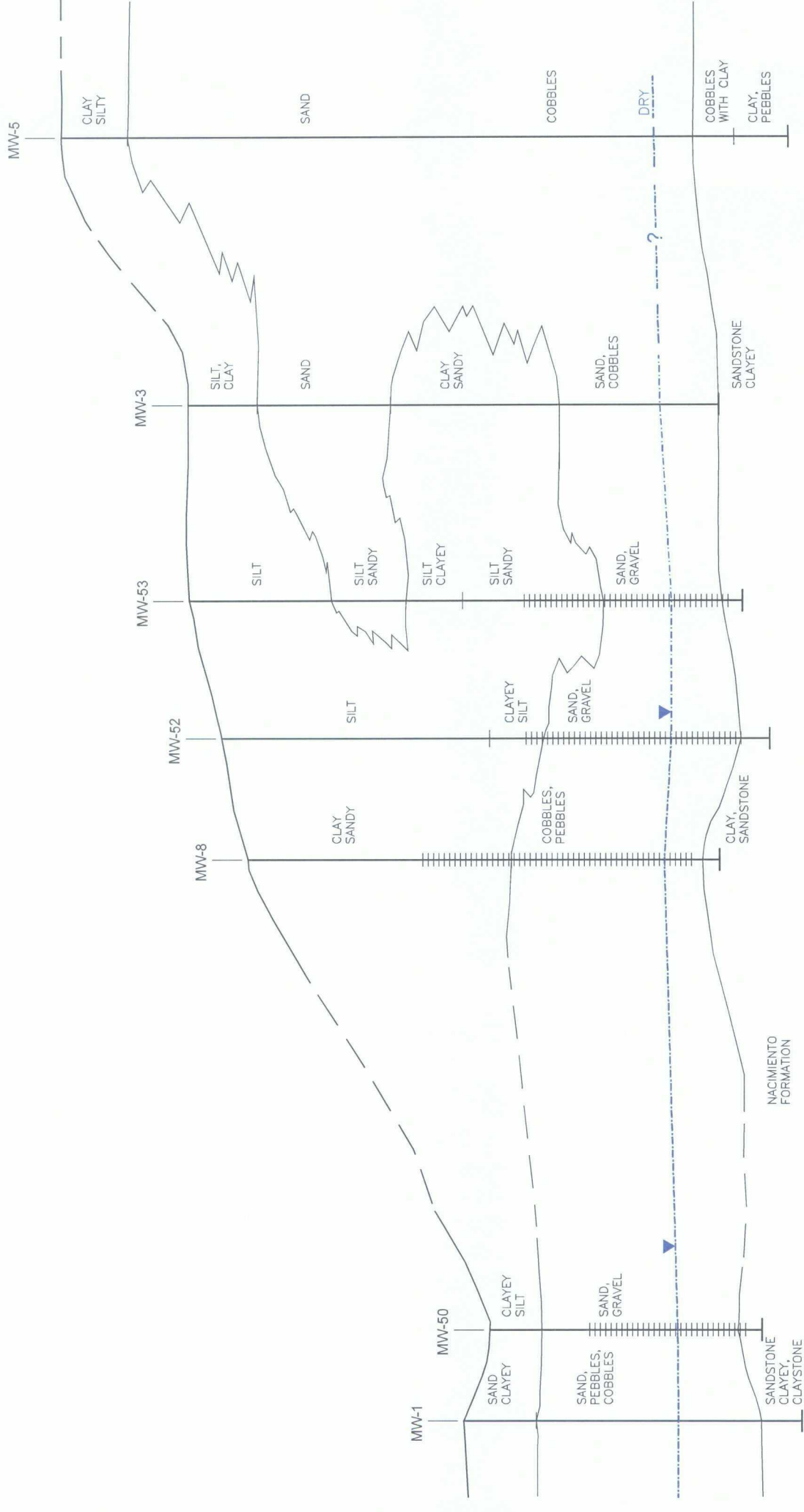
404 Camp Craft Road  
Austin, Texas 78746

SCALE IN FEET  
VERTICAL EXAGGERATION = 20X  
— POTENTIOMETRIC SURFACE MEASURED AUGUST-OCTOBER 2008

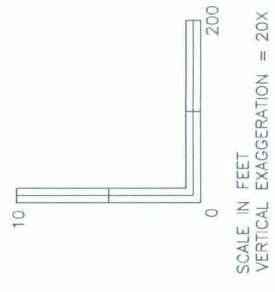
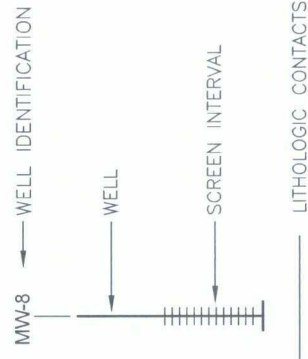
NORTH  
B



SOUTH  
B'



EXPLANATION



PROJ. NO.: Western Refining | DATE: 04/29/09 | FILE: WestRef-B29

FIGURE 4

CROSS SECTION B-B'  
NORTH TO SOUTH  
BLOOMFIELD REFINERY

RPS

404 Camp Craft Road  
Austin, Texas 78746

POTENTIOMETRIC SURFACE MEASURED AUGUST-OCTOBER 2008





|                             |                |                   |
|-----------------------------|----------------|-------------------|
| PROJ. NO.: Western Refining | DATE: 05/21/09 | FILE: WestRef-B19 |
|-----------------------------|----------------|-------------------|

FIGURE 5  
POTENTIOMETRIC SURFACE MAP  
AUGUST 2008  
BLOOMFIELD REFINERY

404 Camp Craft Road  
Austin, Texas 78746

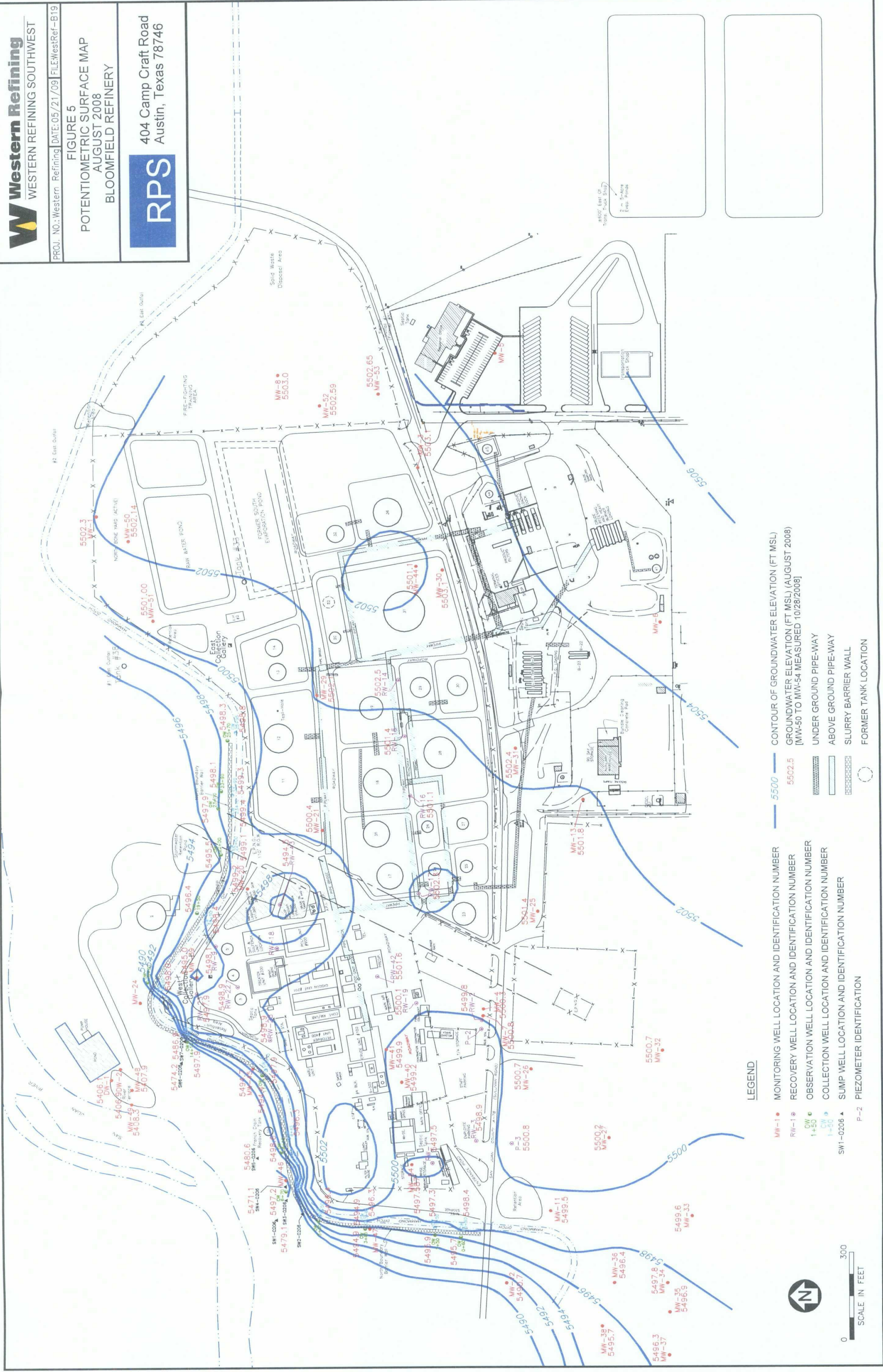


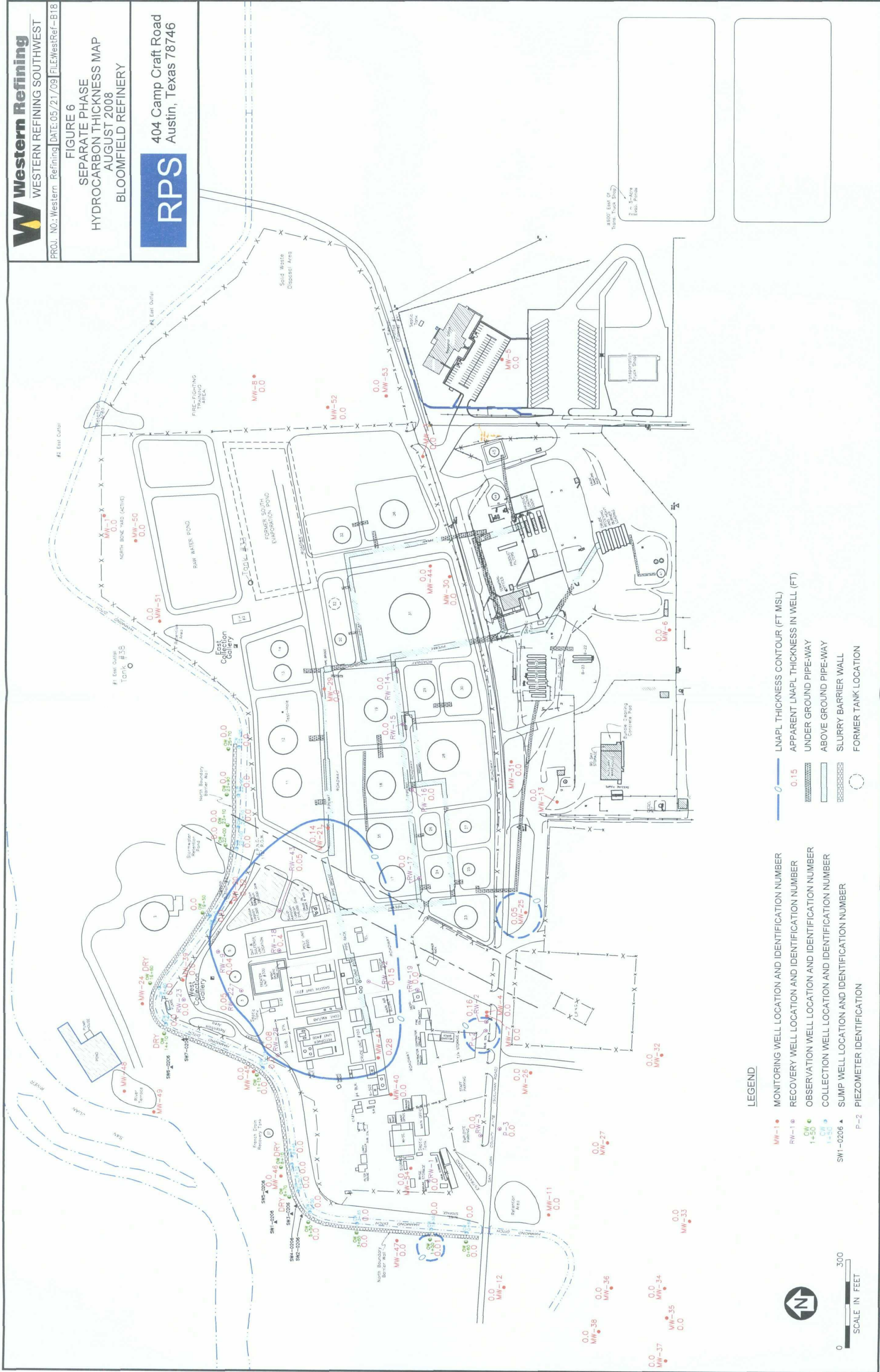


FIGURE 6

SEPARATE PHASE  
HYDROCARBON THICKNESS MAP  
AUGUST 2008  
BLOOMFIELD REFINERY

**RPS**

404 Camp Craft Road  
Austin, Texas 78746







Legend

- Monitoring Well
- Observation Well
- Recovery Well
- Collection Well
- Outfall
- Site
- Approximate Property Line

|          |                  |
|----------|------------------|
| MW-38    | - Well Id        |
| < 0.001  | - Benzene        |
| < 0.001  | - Toluene        |
| < 0.001  | - Ethylbenzene   |
| < 0.0015 | - Xylenes, Total |
| 0.001    | - MTBE           |

Notes:

All concentrations in milligrams per liter (mg/L)

NS = Well is Dry or Not Enough Water to Sample- No sample

NS<sup>2</sup> = Sample Inadvertently not Collected this Sampling Event

NR<sup>1</sup> = No Sample Required - Well Contains Separate Phase Hydrocarbon

NR<sup>2</sup> = No Sample Required per OGD and NMED Conditions

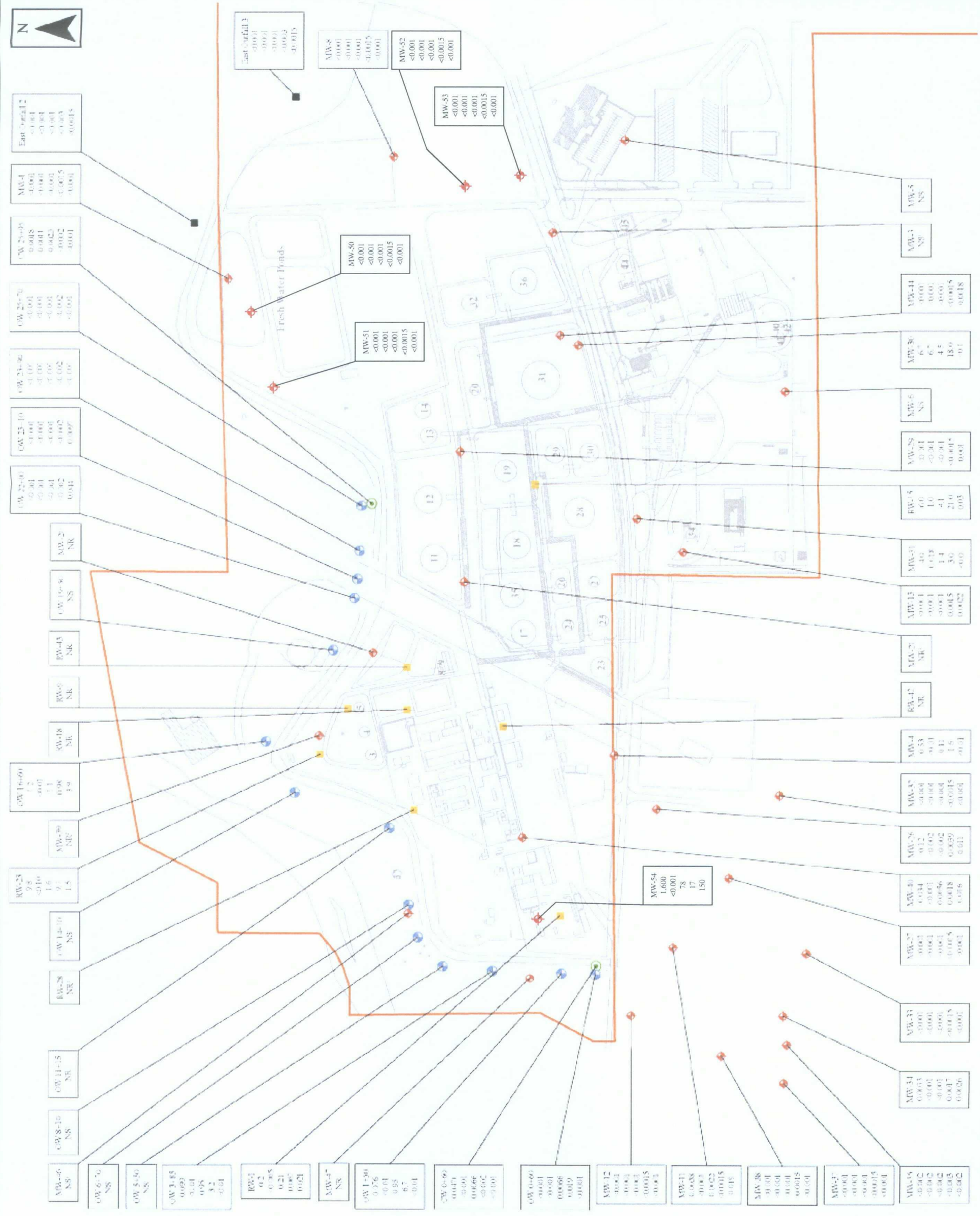


PROJ. NO.: Western Refining DATE: 05/25/09 FILE: WestRef-B32

FIGURE 7

DISSOLVED-PHASE  
GROUNDWATER DATA  
AUGUST AND OCTOBER 2008  
BLOOMFIELD REFINERY

404 Camp Craft Road  
Austin, Texas 78746







Aerial Map Source: Google Map.



0 100  
SCALE IN FEET

LEGEND

- SEEP 1 ♦ SEEP LOCATION AND IDENTIFICATION NUMBER
- MW-24 ● MONITORING WELL LOCATION AND IDENTIFICATION NUMBER
- RW-28 ⊙ RECOVERY WELL LOCATION AND IDENTIFICATION NUMBER
- OW 16+60 ○ OBSERVATION WELL LOCATION AND IDENTIFICATION NUMBER
- CW 16+60 ○ COLLECTION WELL LOCATION AND IDENTIFICATION NUMBER
- SW6-0206 ▲ SUMP WELL LOCATION AND IDENTIFICATION NUMBER

**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 12/10/09 FILE: WestRef-B49

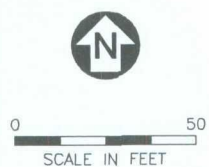
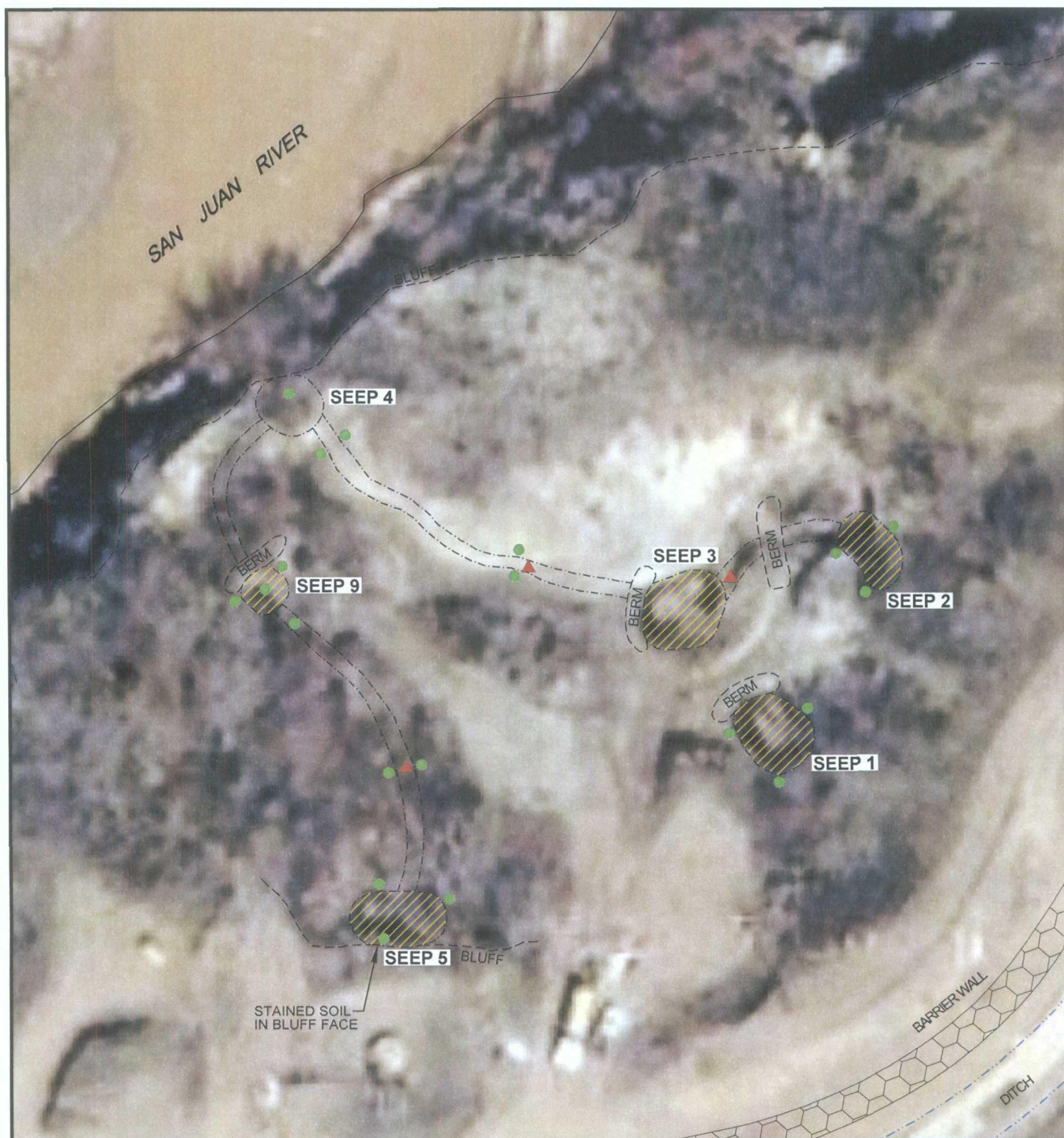
FIGURE 8

AOC Nos. 19, 20 AND 21  
SEEP LOCATION MAP  
BLOOMFIELD REFINERY

**RPS**

404 Camp Craft Road  
Austin, Texas 78746





- LEGEND**
- SEEP 1** SEEP IDENTIFICATION NUMBER
- SURFACE SOIL SAMPLE LOCATION
- ▲ SOIL BORING LOCATION
- ▨ LINED CATCHMENT BASIN
- DRAINAGE PATHWAY

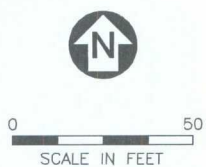
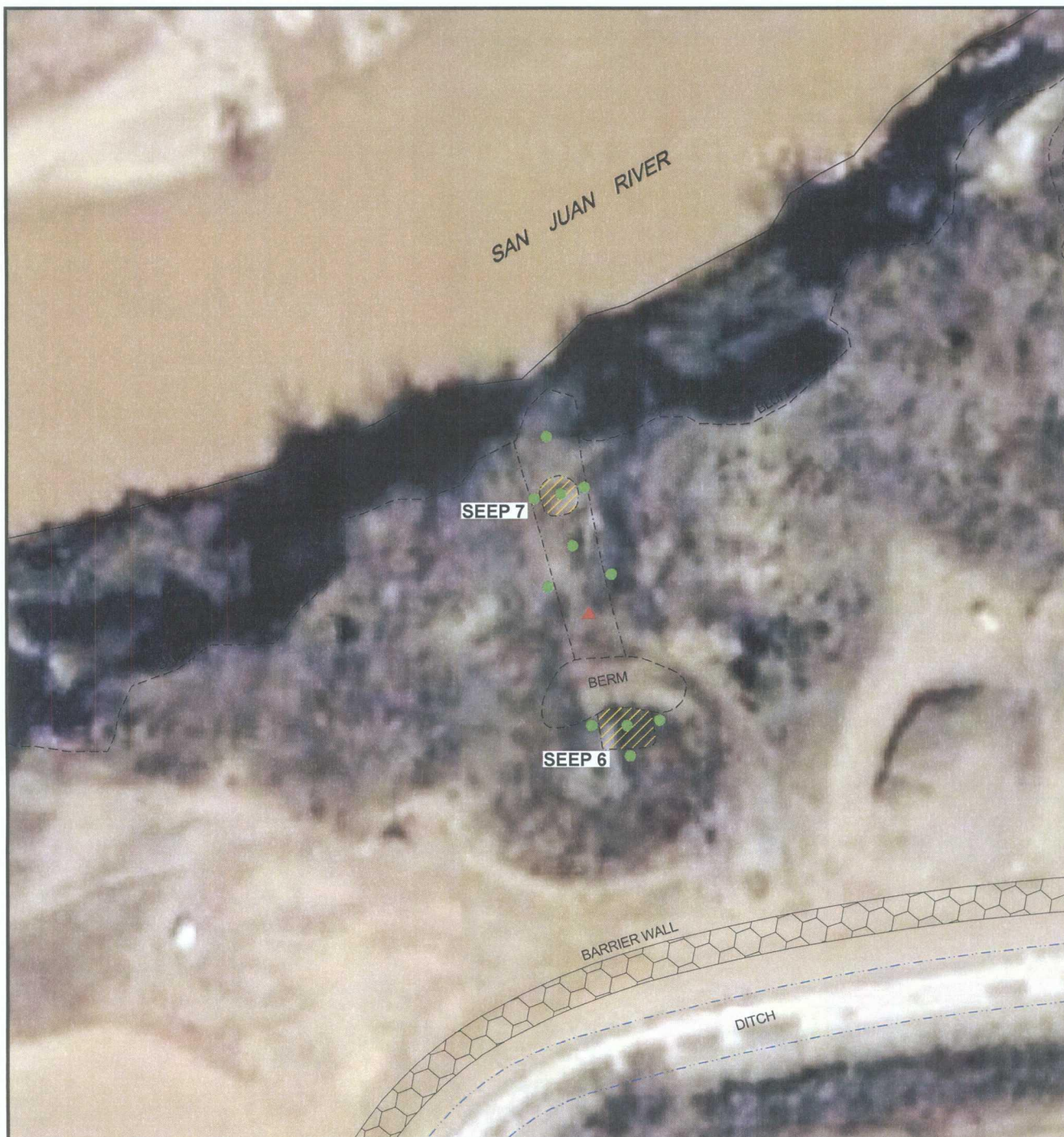
**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining DATE: 12/11/09 FILE: WestRef-A38

FIGURE 9  
AOC No. 19  
SAMPLE LOCATION MAP  
BLOOMFIELD REFINERY

**RPS** 404 Camp Craft Road  
Austin, Texas 78746





- LEGEND**
- SEEP 6** SEEP IDENTIFICATION NUMBER
  - SURFACE SOIL SAMPLE LOCATION
  - ▲ SOIL BORING LOCATION
  - ▨ LINED CATCHMENT BASIN
  - - - DRAINAGE PATHWAY

**Western Refining**  
WESTERN REFINING SOUTHWEST

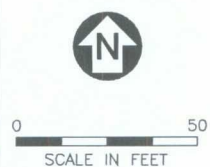
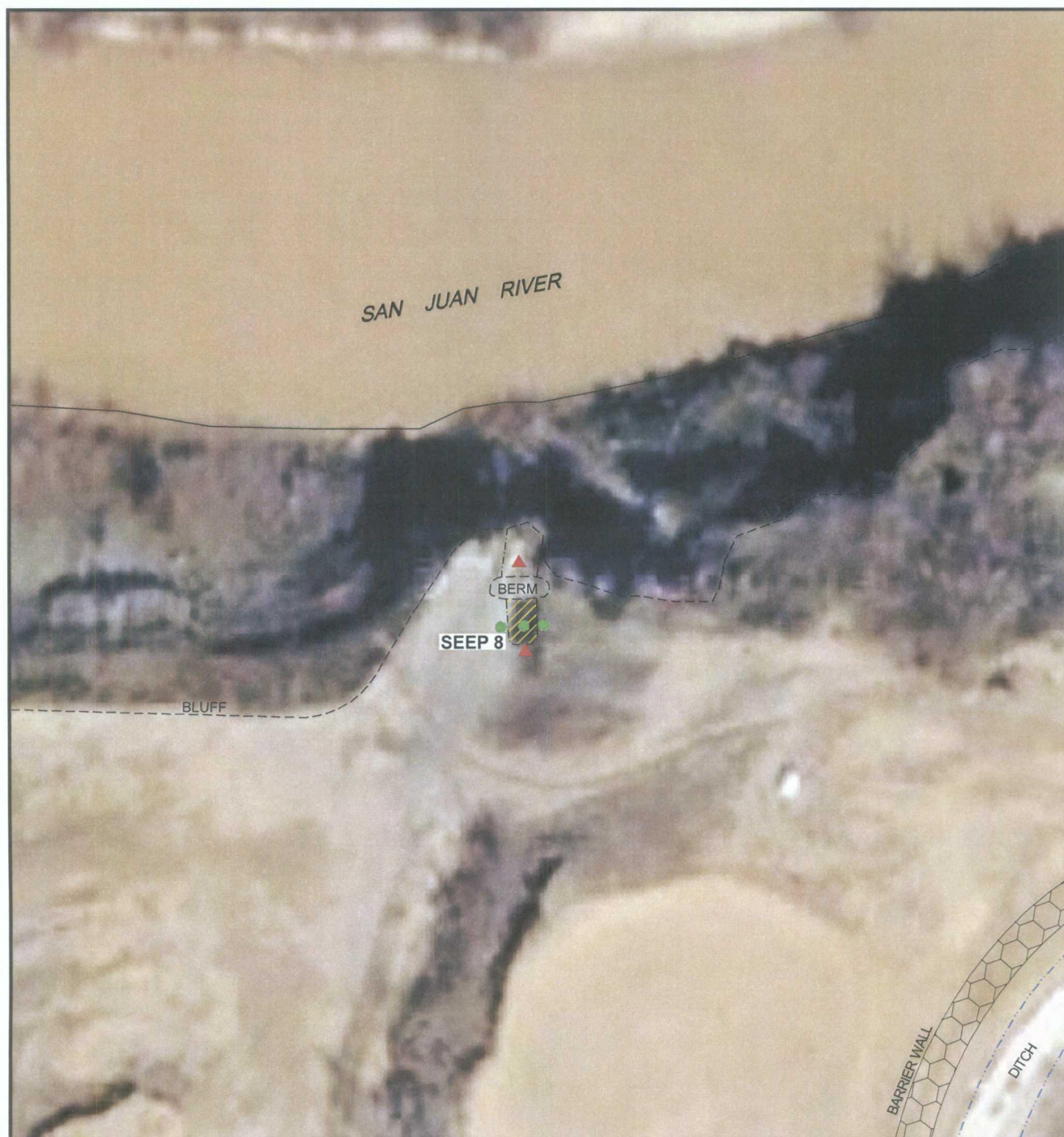
PROJ. NO.: Western Refining | DATE: 12/11/09 | FILE: WestRef-A38

FIGURE 10  
AOC No. 20  
SAMPLE LOCATION MAP  
BLOOMFIELD REFINERY



404 Camp Craft Road  
Austin, Texas 78746





- LEGEND**
- SEEP 8** SEEP IDENTIFICATION NUMBER
- SURFACE SOIL SAMPLE LOCATION
- ▲ SOIL BORING LOCATION
- ▨ LINED CATCHMENT BASIN
- DRAINAGE PATHWAY

**Western Refining**  
WESTERN REFINING SOUTHWEST

PROJ. NO.: Western Refining | DATE: 12/11/09 | FILE: WestRef-A38

FIGURE 11  
AOC No. 21  
SAMPLE LOCATION MAP  
BLOOMFIELD REFINERY



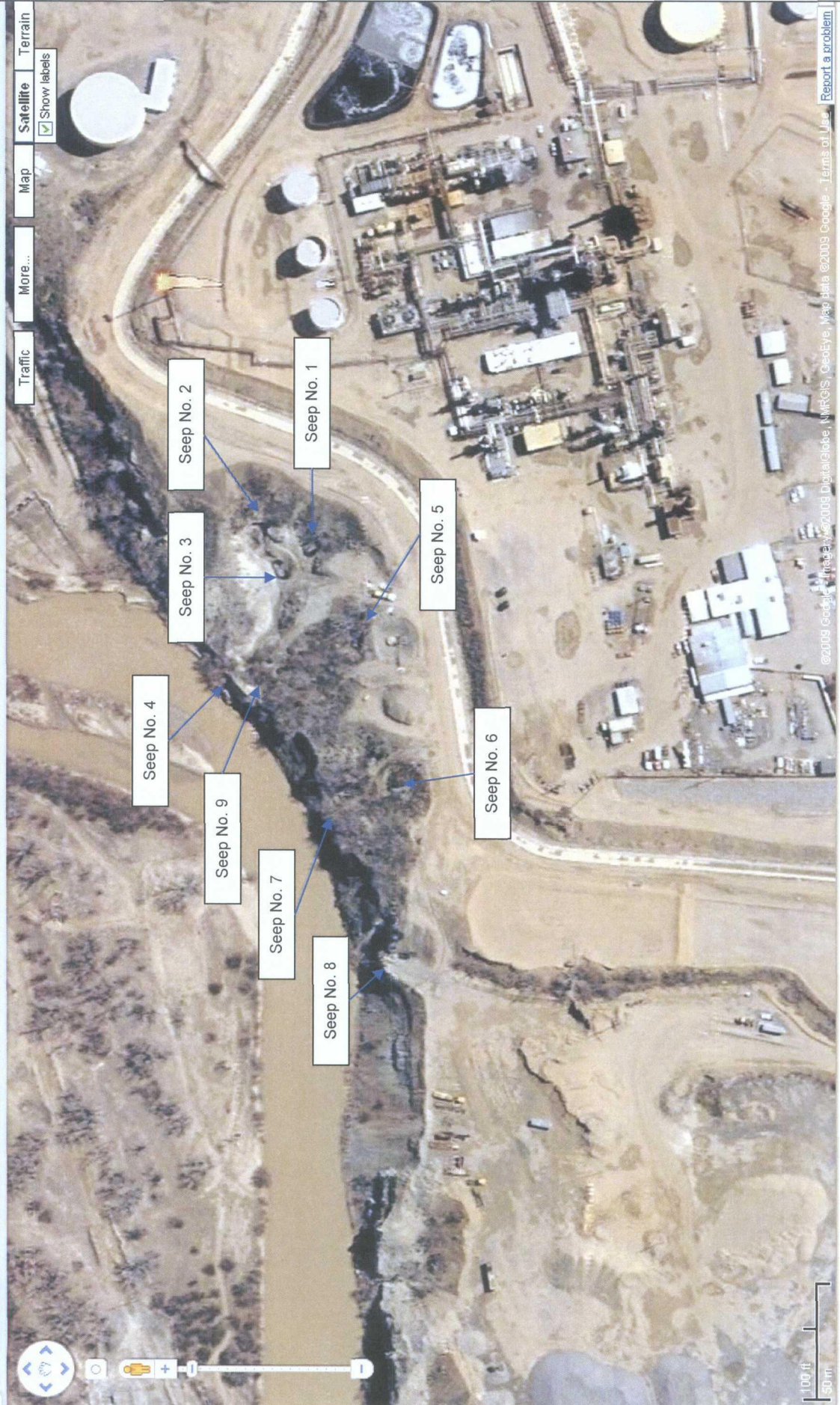
404 Camp Craft Road  
Austin, Texas 78746

# **Appendix A**

---

## **Photographs**

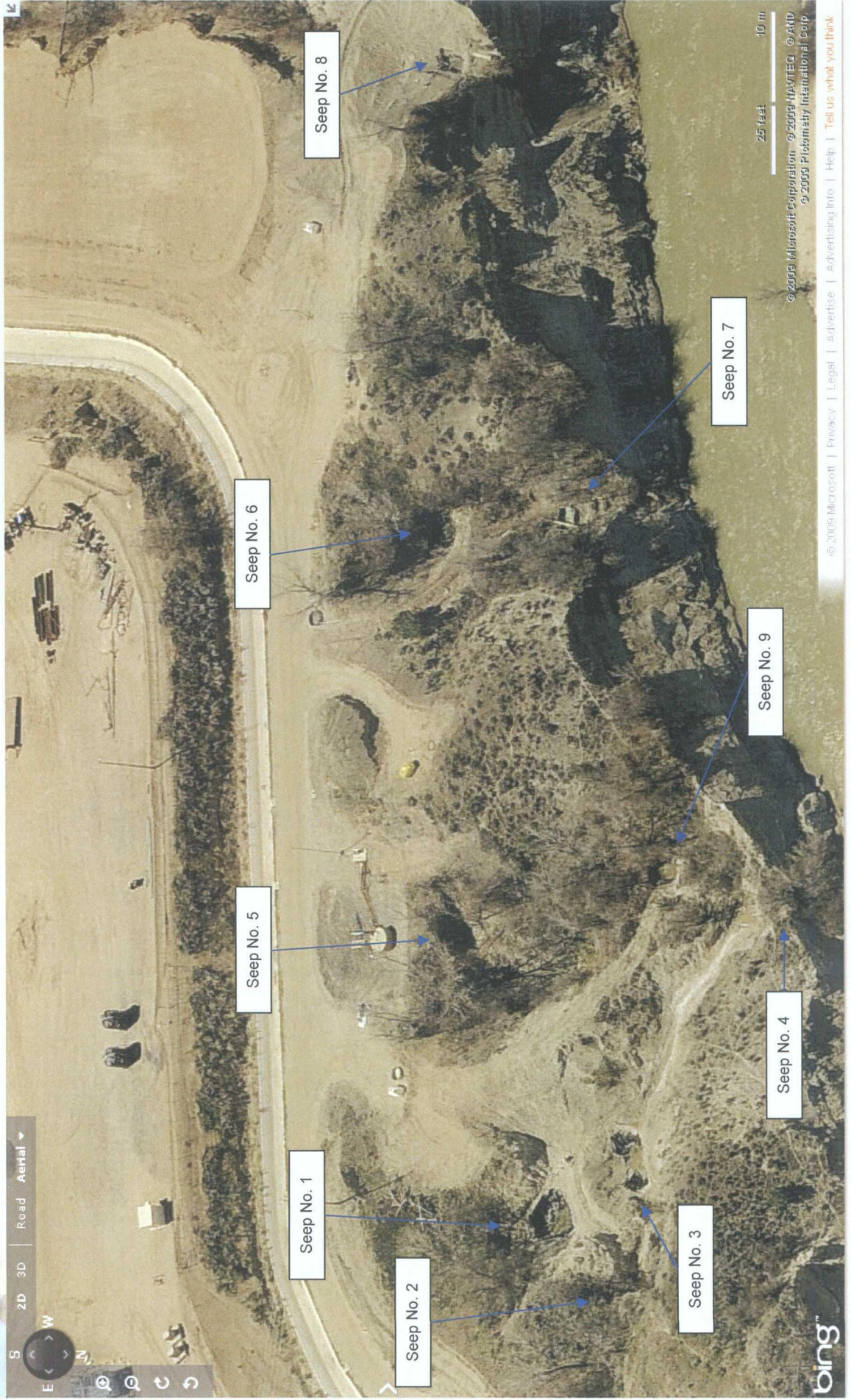




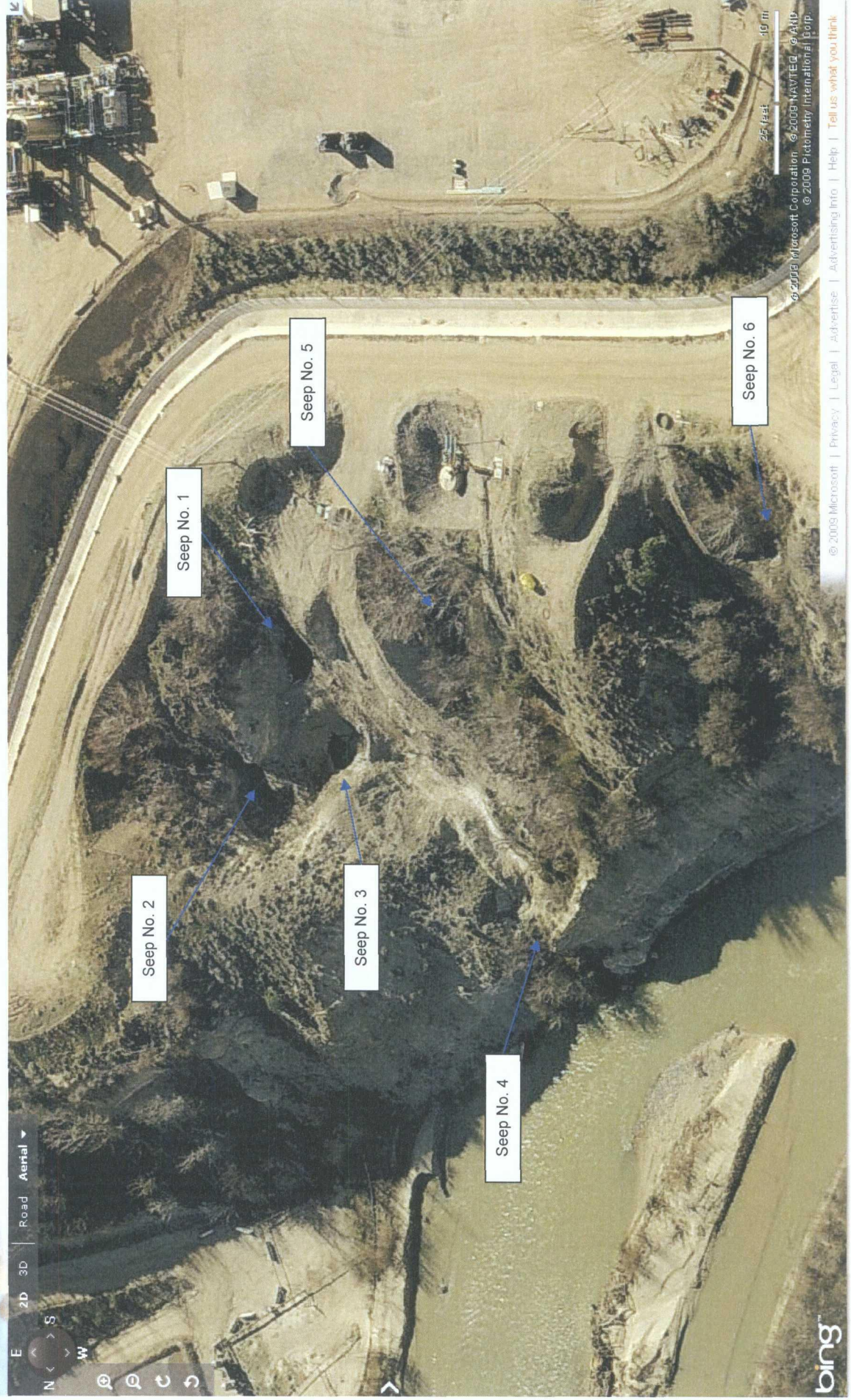
















Bloomfield, NM

bing







AOC No. 19 – Seep and Catchment Basin No. 1



AOC No. 19 – Seep No. 2





AOC No. 19 – Sediment Within Catchment Basin No. 3



AOC No. 19 – Looking Southeast Toward Catchment No. 3 from Catchment No. 4





AOC No. 19 – HDPE Liner of Seep No. 5 Catchment Basin



AOC No. 19 – Stained Soil in Bluff Face of Seep No. 5





AOC No. 19 – Exposed Liner at Seep No. 9



AOC 20 – Seep No. 6 With Water in Catchment Basin No. 6





AOC 20 – Looking North Toward Seep No. 7 From Berm Below Catchment Basin No. 6



AOC 20 – Seep No. 7 and Bluff Overlooking San Juan River





AOC 21 – Seep No. 8 (Below Wood Wall) and Liner of Catchment Basin No. 8



AOC 21 – Looking North Over San Juan River Bluff From Catchment Basin No. 8

## **Appendix B**

---

### **Investigation Derived Waste (IDW) Management Plan**



## **IDW Management Plan**

All IDW will be properly characterized and disposed of in accordance with all federal, State, and local rules and regulations for storage, labeling, handling, transport, and disposal of waste. If soils are found to be non-hazardous and concentrations of constituents are less than the NMED residential soil screening levels, then soils may be reused on-site pursuant to the approval of the NMED. The IDW may be characterized for disposal based on the known or suspected contaminants potentially present in the waste. It is assumed that there are no listed wastes present in environmental media at any of the planned investigation areas.

A dedicated decontamination facility will be setup prior to any sample collection activities. The decontamination facility will be designed so as to capture and contain all decontamination fluids (e.g., wash water and rinse water) and foreign materials washed off the sampling equipment. The fluids will be pumped directly into suitable storage containers (e.g., labeled 55-gallon drums), which will be located at satellite accumulation areas until the fluids are disposed in the refinery wastewater treatment system upstream of the API separator. The solids captured in the decontamination facility will be shoveled into 55-gallon drums and stored at the designated satellite accumulation area pending proper waste characterization for off-site disposal.

Any solids generated (e.g., used soil cores) will be characterized by testing to determine if there are any hazardous characteristics in accordance with 40 Code of Federal Regulations (CFR) Part 261. This includes tests for ignitability, corrosivity, reactivity, and toxicity. If the materials are not characteristically hazardous, then further testing will be performed pursuant to the requirements of the facility to which the materials will be transported. Depending upon the results of analyses for individual investigation soil samples, additional analyses may include TPH and polynuclear aromatic hydrocarbons. All miscellaneous waste materials (e.g., discarded gloves, packing materials, etc.) will be placed into the refinery's solid waste storage containers for off-site disposal.