# **GW - 28**

2019 AGWMR (1)

2020

From: <u>Dade</u>, Randy

To: <u>Chavez, Carl J, EMNRD</u>

Cc: <u>Dade, Randy</u>

Subject: [EXT] 2019 Annual Discharge Report, HollyFrontier Navajo Refining LLC, Artesia Refinery, Discharge Permit GW-

28

**Date:** Tuesday, June 16, 2020 7:12:44 AM

Attachments: 2019 Annual Discharge Report Transmittal Letter.pdf

2019 GW-028 Annual Discharge Report FINAL 06152020.pdf

#### Carl,

Please find attached the 2019 Annual Discharge Report, HollyFrontier Navajo Refining LLC, Artesia Refinery, Discharge Permit GW-28. I will be uploading an electronic version to the OCD website. If you have any comments or questions, please contact me.

Thanks for all your help that you provide, Randy.

Lewis R. (Randy) Dade Environmental Specialist The HollyFrontier Companies 501 E. Main / P.O. Box 159 Artesia, NM 88210 / 88211-0159 575-746-5281 (o) 575-703-4735 (c) 575-746-5451 (f)

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June 15, 2020

Mr. Carl Chavez
New Mexico Energy, Minerals and Natural Resources Department
Oil Conservation Division
5200 Oakland Avenue N.E., Suite 100
Albuquerque, NM 87113

Re: Submittal of the 2019 Annual Discharge Report and the 2019 Annual Groundwater Monitoring Report for the HollyFrontier Navajo Refining LLC, Artesia Refinery Discharge Permit GW-028

Dear Mr. Chavez:

Please find attached the 2019 Annual Discharge Report and the 2019 Annual Groundwater Monitoring Report, which fulfill requirements of Section 2.E of Discharge Permit GW-028. No hard copy will be submitted at this time.

If you have any questions or comments regarding this report, please feel free to contact me at 575-746-5487 or Robert Combs at 575-746-5382.

Sincerely,

Scott M. Denton

Environmental Manager

HollyFrontier Navajo Refining LLC

cc: HollyFrontier: R. Combs, J. Leik, R. Dade TRC: J. Speer, C. Smith, D. Helbert



# 2019 Annual Discharge Permit Report, GW-028

June 15, 2020

# **HollyFrontier Navajo Refining LLC Artesia Refinery, GW-028**

#### **Prepared For:**

HollyFrontier Navajo Refining LLC 501 E Main Street, Artesia, NM 88210

#### **Prepared By:**

TRC 505 East Huntland Drive, Suite 250 Austin, TX 78752







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#### **ATTACHMENTS**

Attachment A 2019 Annual Groundwater Monitoring Report, February 2020 (Separate Electronic File)



#### ABBREVIATION AND ACRONYM LIST

**ACO** Agreed Compliance Order

AOC Area of concern

**bbl** barrel

**bpd** barrels per day

BTEX Benzene, toluene, ethylbenzene, and xylene

**CGWSL** Critical Groundwater Screening Level

COCs Constituents of concern
DRO Diesel range organics
gpm Gallons per minute

**GRO** Gasoline range organics

**HFNR** HollyFrontier Navajo Refining LLC

MTBE Methyl tert-butyl ether

NMAC New Mexico Administrative Code
NMED New Mexico Environment Department

O&M Operation and maintenance
OCD Oil Conservation Division
ORP Oxidation-reduction potential
PSH Phase-separated hydrocarbon

**PCC** Post Closure Care

**POTW** Publicly Owned Treatment Works

**Praxair** Praxair Services, Inc.

**RCRA** Resource Conservation and Recovery Act

RO Reverse osmosis

**SPLP** Synthetic precipitation leaching procedure

**SSL** Soil screening levels

SVOC Semi-volatile organic compound SWMU Solid Waste Management Unit

TDS Total dissolved solids
TEL Tetra Ethyl Lead

TPH Total petroleum hydrocarbons
UIC Underground Injection Control
VOC Volatile organic compound

WDW Water Disposal Well

WQCC Water Quality Control Commission



#### Introduction

This report was prepared to fulfill the requirement in Section 2.E. of the Discharge Permit GW-028 (GW-028) for the HollyFrontier Navajo Refining LLC (HFNR) Artesia Refinery (refinery) located at 501 East Main Street in Artesia, New Mexico. The requirement specifies that an Annual Report be submitted to the Oil Conservation Division (OCD) by June 15 following the reporting (calendar) year and should include:

- 1. Summary of major refinery activities and events.
- 2. Summary of all discharge activities.
- 3. Summary of all leaks, spills, and releases and corrective actions taken.
- 4. Summary of discovery of any new vadose zone or groundwater contamination.
- 5. Summary of wastewater volumes disposed of, sold, or treated onsite.
- Documentation regarding the closure of any Underground Injection Control (UIC) Class V wells.
- 7. A description of groundwater monitoring and remediation activities conducted throughout the year.
- 8. Summary tables of groundwater data.
- 9. Copies of laboratory analytical data sheets with quality assurance/quality control information.
- 10. Contour maps for each aquifer depicting the potentiometric gradient for each monitoring event.
- 11. Isoconcentration maps of major constituents of concerns (COCs) for each monitoring event.
- 12. Phase-separated hydrocarbon (PSH) thickness isopleth maps for each monitoring event.
- 13. Plots of static water elevation versus time in key wells.
- 14. Tabulation of the volumes of PSH removed.
- 15. Conclusions and recommendations.



#### 1.0 Major Refinery Activities for 2019

The refinery conducted normal operations during 2019. Additional capital projects were completed to improve operability. No new tanks or refinery units were built in 2019.

#### 1.1 Discharge Permit GW-028 Modifications

The previous GW-028 (dated August 22, 2012) was set to expire on October 21, 2016. HFNR submitted an application for renewal of and modification to GW-028 on June 23, 2016 (at least 120 days prior to expiration). OCD notified HFNR that the application was administratively complete on July 28, 2016, and HFNR proceeded to complete all required public notices.

On September 9, 2016, OCD notified HFNR that the renewal application did not propose a definitive alternative, or the information required to evaluate such alternative, to replace land application of reverse osmosis (RO) reject water discharge. As such, OCD would not issue an approval or disapproval of the renewal application until such information was provided. On September 23, 2016, HFNR entered into Agreed Compliance Order (ACO) No. WQA-OCD-CO-2016-1 (the 2016 Order) which allowed continued RO discharge operations (per Condition III.1.a.iii of the 2016 Order) while progressing with the Discharge Permit renewal application.

On October 21, 2016, HFNR notified OCD of the selection of underground injection as the alternative disposal method for the RO reject stream in accordance with Condition III.1.a.i of the 2016 Order. HFNR submitted a revised Discharge Permit renewal application reflecting the selection of underground injection (through a Class I disposal well) as the alternative disposal method on January 13, 2017. OCD issued a renewal to GW-028 on May 25, 2017, the Order was terminated on June 21, 2017, and a modification of GW-028 was issued on June 29, 2017. The renewed GW-028 included a stipulation that land application of the RO reject water must cease upon the completion of the new Class I disposal well, but not later than October 31, 2018. The OCD issued modifications of GW-028 on October 25, 2018, and December 14, 2018, which extended the deadline for land application of RO reject water due to delays in operational completion of the new Class I disposal well.

HFNR is continuing to work with OCD for permitting the sale of treated refinery effluent for reuse in the oilfield.

#### 1.2 RO Reject Water Discharge

From January 1 to January 22, 2019, HFNR discharged RO reject water to the refinery's onsite fields, in accordance with GW-028 which allows for discharge of RO reject water to the refinery's onsite fields. The land application of RO reject water was discontinued after January 22, 2019, as described below. GW-028 requires sampling and analyzing RO reject water for Water Quality Control Commission (WQCC) constituents. HFNR collected a grab sample of RO reject water from the point of discharge on January 18, 2019, prior to the cessation of land application of RO reject water. Laboratory analytical reports are provided in Appendix A.3.



#### 1.3 Injection Well WDW-4 (Class I Disposal Well)

HFNR selected to install a fourth injection well (WDW-4) as an alternate disposal method for the RO reject water, as specified in the revised permit application. Installation of WDW-4 and associated piping was completed in late 2018 and the well became operational on January 16, 2019. The well was utilized for injection of water beginning on January 23, 2019.

#### 1.4 RO Reject Fields Investigation and Abatement Plan

On August 20, 2015, HFNR submitted a *Reverse Osmosis Reject Fields Hydrogeologic and Water Quality Evaluation* memo to the OCD that fulfilled the Site investigation requirements of Section 6.D of the former GW-028 (dated August 22, 2012). A subsequent revision to this memo was submitted to OCD on January 19, 2016, to provide corrections to the RO reject stream water quality results. HFNR met with the OCD and New Mexico Environment Department (NMED) at the OCD office on March 11, 2016, to discuss the results of the background groundwater evaluation (submitted to NMED and OCD in September 2015) as well as the hydrogeologic model and loading report. No agreement was reached regarding the results and recommendations of either of these evaluations.

HFNR began discussions with OCD in March 2017 regarding the potential to abate WQCC constituents in the RO reject water and in the RO reject discharge fields via phytoremediation. HFNR conducted a phytoremediation feasibility study at the RO reject fields from August 2017 to March 2018. Results of the phytoremediation feasibility study were documented in the *Phytoremediation Feasibility Study Summary Report* that was included as an appendix to the required Abatement Plan, described below.

GW-028 requires discharge of RO reject water to the fields to cease upon the completion of a Class I injection well and submittal of an Abatement Plan within 60 days of cessation of discharge of RO reject water to the fields. As described above, WDW-4 became operational in January 2019 and land application of RO reject water ceased after January 22, 2019.

HFNR submitted to OCD a Stage 1 Abatement Plan for the Reverse Osmosis Reject Discharge Fields on March 21, 2019 and an Amendment of the March 2019 Stage 1 Abatement Plan for the Reverse Osmosis Reject Discharge Fields on May 24, 2019. HFNR is characterizing the RO reject fields after cessation of land application of RO reject water to support development of a Stage 2 Abatement Plan, as agreed by OCD and HFNR in a meeting on May 16, 2019, and in accordance with the May 2019 Amendment of the March 2019 Stage 1 Abatement Plan for the Reverse Osmosis Reject Discharge Fields.

#### 2.0 Summary of Discharge Activities

HFNR's primary discharges are treated wastewater from the wastewater treatment plant (WWTP) effluent and the RO reject water. The WWTP effluent is discharged to HFNR's injection wells (WDW-1, WDW-2, WDW-3 and WDW-4) and to the City of Artesia's Publicly Owned Treatment Works (POTW). RO reject water was land applied from January 1 to January 22,



2019. After January 22, 2019, RO reject water is further processed in a secondary RO unit, which produces a permeate stream which is utilized in the refinery's cooling towers, and the reject stream from that unit is ultimately discharged to HFNR's injection wells or the City of Artesia's POTW. The details of each discharge are provided in the following sub-sections.

#### 2.1 Injection Wells

The injection rates, volume, and quality of treated wastewater disposed of in the injection wells are reported quarterly to OCD, in addition to monthly C-115 reports. Injection rates and volumes are also summarized in a table provided as Appendix A.1. The total injected water volume for 2019 was 6,240,894 barrels.

#### **2.2 POTW**

The flow rates and volumes of treated wastewater discharged to the City of Artesia POTW are recorded monthly and provided as Appendix A.2. The total transferred water volume for 2019 was 1,156,176 gallons or 27,528 barrels.

HFNR continued to discharge the blow-down from cooling towers to the City of Artesia POTW in 2019. The total volume of blow-down discharged to the City of Artesia POTW based on an average rate of 86 gallons per minute (gpm) is estimated to be 45,800,640 gallons, or 1,090,491 barrels.

#### 2.3 Reverse Osmosis Reject Water

The RO reject water was land applied under GW-028 to onsite fields from January 1, 2019 through January 22, 2019, when the new injection well, WDW-4, was utilized for RO reject water disposal. The supply lines to the RO fields were capped and the main valve was shut off, preventing further discharge of RO reject water to the fields.

The RO process is fed by fresh groundwater provided by either the refinery's agricultural supply wells or purchased from the City of Artesia. The RO reject fluids contain concentrated salts (primarily chloride, fluoride, and sulfate) and elevated total dissolved solids (TDS). The stream was sampled on January 18, 2019 in accordance with GW-028. Laboratory analytical reports are provided in Appendix A.3.

The RO reject fluid flow rate was continuously recorded with the process historian and copies are provided in Appendix A.3. Based on the data from the process historian and on the logs, the total discharged RO reject water volume for the 22 days of land discharge in January 2019 was 10,289,760 gallons, or 244,994 barrels. The average daily discharge rate was 11,136 barrels per day. There were no exceedances of the permitted discharge rate in 2019.

#### 3.0 Summary of All Leaks, Spills, and Releases

The refinery had four reportable spills under GW-028 in 2019. Each spill was reported to the OCD and addressed as described below.



#### 3.1 March 26, 2019 - Tank 106 Release

Approximately 40 to 50 barrels of sour water was released on March 26, 2019, from Tank 106 due to a local gauge and transmitter malfunction. The release was entirely contained within the Tank 106 secondary containment, which consists of earthen berms. The tank level was reduced to prevent further release. Over 40 barrels of free liquids were recovered and placed into the refinery wastewater treatment system, upstream of the oil/water separator. Impacted soil was removed and placed into roll off bins. The initial C-141 Release Notification for this release was submitted to OCD on March 27, 2019.

Soil assessment activities were conducted in June 2019. Assessment results indicated benzene, toluene, ethylbenzene, and xylene (BTEX), chloride, total petroleum hydrocarbons (TPH), and benzene were present in soil at concentrations above their respective closure criteria, but below their applicable NMED Construction Worker soil screening levels (SSLs). TPH was present in soil above NMED Construction Worker SSLs. Worker protection corrective measures, including placement of clean soil and gravel over the release area, were implemented. The impacts were located in an area that contains sensitive refinery equipment and that is already identified as area of concern 3 (AOC 3) in the refinery's RCRA Post-Closure Care Permit (PCC) Permit and is therefore already subject to investigation and corrective action under the direction of the NMED. Therefore, HFNR requested a variance from the requirements to assess and remediate to 19.15.29.12 NMAC Table 1 standards. A Site Characterization, Assessment, and Closure Report was submitted to the New Mexico OCD on September 20, 2019, and included a request for a variance to 19.15.29.11(A)(5), 19.15.29.11(B), and 19.15.29.12 NMAC and a final C-141 (Site Assessment/Characterization and Closure).

#### 3.2 May 28, 2019 – Cooling Tower Blowdown Sampling Station Release

Non-hazardous cooling tower blowdown water was released on May 28, 2019, from plastic tubing near a sampling station on the cooling tower blowdown line. The line was shut down and the tubing was repaired. The initial C-141 Release Notification for this release was submitted to New Mexico OCD on May 30, 2019.

Soil assessment activities were conducted in June 2019. Assessment results indicated BTEX, fluoride, sulfate, phenol, and arsenic concentrations were below their respective closure or screening criteria. Assessment results indicated chloride and TPH were present in soil at concentrations above their respective closure criteria, but significantly below their applicable NMED Construction Worker SSLs. The impacts were located in an area containing sensitive refinery infrastructure and distribution, and variability of TPH and chloride concentrations across the release are indicate they are not attributable to the May 2019 release. Therefore, HFNR requested that corrective action of impacted soil be deferred until the infrastructure is removed in accordance with 19.15.29.12 NMAC. A Site Characterization, Assessment, and Closure Report was submitted to the New Mexico OCD on August 28, 2019, and included the request for deferred corrective action and a final C-141 (Site Assessment/Characterization and Closure).

During a call in December 2019, OCD confirmed that 19.15.29 NMAC is not applicable to the refinery. During that call, it was determined that in order to address refinery releases, HFNR



must review 20.6.2 NMAC to determine a basis for OCD to allow releases to remain in place or accept a risk-based approach to remediation after release characterization. Additionally, HFNR must develop environmental investigation, characterization, and remediation guidelines for OCD to consider and approve. A Draft *HFNR Release Response and Characterization Plan* was submitted to the OCD on June 1, 2020 for OCD's review and comment. HFNR will continue to monitor shallow groundwater immediately beneath the release area on a semi-annual basis as part of the refinery's groundwater monitoring program.

#### 3.3 September 3, 2019 – Cooling Tower Blowdown Sampling Station Release

Non-hazardous cooling tower blowdown water was released on September 3, 2019, from tubing near a sample station on the refinery's cooling tower blowdown line to the City of Artesia's POTW. The release location and impacted area are nearly identical to the May 28, 2019 cooling tower blowdown sampling station release, and extent of the release area was entirely contained within the refinery fence line. The sample station was isolated, the tubing was repaired, and the impacted area was marked and defined. The initial C-141 Release Notification for this release was submitted to New Mexico OCD on September 4, 2019. A final C-141 was submitted on September 9, 2019.

Further action for this release will be consistent with the May 2019 cooling tower release. HFNR will continue to monitor shallow groundwater immediately beneath the release area on a semi-annual basis as part of the refinery's groundwater monitoring program.

#### 3.4 November 8, 2019 - Tank 401

During a tank inspection by Praxair Services, Inc. (Praxair), moist soil was observed at the base of T-401, an above ground storage tank containing gasoline blendstock. Praxair tested existing probes for detection of their inoculant. Initial tests suggested that the inoculant detections may have been only from the interstitial space between the two tank floors. Subsequent testing indicated tracer compound was detected below the secondary floor at probe 6, as described in the Praxair Tracer Tight Leak Test Report, dated November 8, 2019, submitted with the initial C-141 Release Notification, submitted to the New Mexico OCD on November 22, 2019. The release was not confirmed to be greater than 5 barrels. The tank was removed from service and was emptied for inspection. A final C-141 report is in development. HFNR will continue to monitor shallow groundwater immediately beneath the release area on a semi-annual basis as part of the refinery's groundwater monitoring program.

#### 4.0 Summary of New Groundwater Contamination

Groundwater contamination and changes in existing constituents are discussed in Section 7 of the 2019 Annual Groundwater Monitoring Report that was submitted to the NMED on February 28, 2020 (and attached to this report). Groundwater conditions measured during 2019 semiannual events were generally consistent with historical results as summarized below:

 The presence and distribution of PSH were generally consistent with previous monitoring results, with minor fluctuations. PSH thicknesses across the refinery are stable to



declining over time with the exception in select wells (KWB-10R, MW-112, MW-127, and MW-128, MW-137, MW-138) located in the Field East of Refinery and North Refinery that are attributed to reductions in groundwater elevations. PSH thicknesses are inversely affected by fluctuations in groundwater elevations, which generally decreased 2017 through 2019, consistent with drought conditions across New Mexico during this time.

- Concentrations of COCs in groundwater have generally remained stable over time, although increasing trends were noted in select wells in specific areas of interest. The limited number of increasing COC concentration trends observed since 2011 have generally exhibited stabilizing trends over the most recent sampling events. During 2019 and previous years, the following COCs were detected in groundwater at concentrations in exceedance of their respective critical groundwater screening level (CGWSL):
  - o TPH gasoline range organics (GRO) and diesel range organics (DRO);
  - Select volatile organic compounds (VOCs) including target COCs benzene,
     toluene, ethylbenzene, xylenes, methyl tert-butyl ether (MTBE), and naphthalene;
  - Select total metals including target COC arsenic; and
  - o Water quality parameters chloride, fluoride, sulfate, TDS, and nitrate/nitrite.
- Semi-volatile organic compounds (SVOCs) were detected in groundwater at concentrations in exceedance of their respective CGWLS in samples collected from select wells in the vicinity of Solid Waste Management Units (SWMUs) 20 and 22 (North Refinery and TEL wells).
- Many of the concentrations of inorganic COCs (manganese, chloride, fluoride, nitrate/nitrite, sulfate, and TDS) noted as "exceedances" of CGWSLs in 2019 may actually be similar to and reflective of background groundwater concentrations, as detailed in the background groundwater evaluation that was submitted to NMED and OCD in September 2015.
- The PSH and groundwater recovery system operated throughout 2019; more information is provided in Section 14 below.

## 5.0 Summary of All Wastewater Volumes Disposed of, Sold, or Treated Onsite

No waste is disposed, sold, or treated onsite.

As described above, wastewater is treated in the refinery WWTP and discharged to either HFNR's injection wells (WDW-1, WDW-2, WDW-3, or WDW-4) or the City of Artesia POTW, both of which are located outside of the refinery (i.e., offsite). The onsite WWTP treated approximately 6,268,422 barrels of wastewater in 2019.

As described above, RO reject water was land applied to onsite fields from January 1 to January 22, 2019. Approximately 244,994 barrels of RO reject water were applied to the RO fields in 2019. RO reject water is no longer applied to onsite fields and is now utilized in the



refinery's cooling towers and ultimately discharged to HFNR's injection wells or the City of Artesia's POTW.

#### 6.0 Documentation Regarding the Closure of Any UIC Class V Wells

No UIC Class V wells were closed during 2019.

# 7.0 A Description of Groundwater Monitoring and Remediation Activities Conducted Throughout the Year

Groundwater monitoring and remediation activities conducted at the refinery in 2019 are described in the attached 2019 Annual Groundwater Monitoring Report. Groundwater monitoring activities, including sample collection procedures, decontamination procedures, sample handling procedures, and investigation-derived waste management, are described in Section 2 of the 2019 Annual Groundwater Monitoring Report. Remediation activities, including PSH recovery, are described in Section 6 of the 2019 Annual Groundwater Monitoring Report.

#### 8.0 Summary Tables of Groundwater Data

Summary tables of groundwater data including water quality, purging parameters, groundwater elevation, and PSH thickness are provided in the attached *2019 Annual Groundwater Monitoring Report*, as specified below.

#### 8.1 Well Gauging Results (Groundwater Elevation and PSH Thickness)

Well gauging results for both 2019 semiannual monitoring events are presented in Table 1 of the attached 2019 Annual Groundwater Monitoring Report. Well gauging results include depth to water measurements, depth to PSH (if present) measurements, and groundwater elevations. Well gauging results for routine PSH recovery operation and maintenance (O&M) activities are summarized in Appendix F of the 2019 Annual Groundwater Monitoring Report.

#### 8.2 Field-Measured Purging Parameters

Groundwater quality parameters measured in the field at each well during 2019 groundwater purging and sampling activities are summarized in Table 2 of the attached 2019 Annual Groundwater Monitoring Report. Groundwater quality parameters include pH, temperature, specific conductance, oxidation-reduction potential (ORP), dissolved oxygen, and turbidity. Observations of relative water quality (color and odor) are also included in Table 2 of the 2019 Annual Groundwater Monitoring Report.



#### 8.3 Laboratory Analytical Results (Water Quality)

Laboratory analytical results of all wells sampled in 2019, and during at least the three previous sampling events, are summarized in Tables 4A through 4D of the attached *2019 Annual Groundwater Monitoring Report* as follows:

- Table 4A Total petroleum hydrocarbons (GRO and DRO) and select VOCs (VOCs that have had at least one detected value reported above the CGWSL in more than one well in 2019)
- Table 4B Total Metals
- Table 4C Water quality parameters (TDS, nitrate/nitrite, major cations, major anions) and Cyanide
- Table 4D Select SVOCs (SVOCs that have had at least one detected value in at least one well in 2019)

Analytical results of all detected COCs are summarized in tables that are included in Appendix C of the attached 2019 Annual Groundwater Monitoring Report.

# 9.0 Copies of Laboratory Analytical Data Sheets with Quality Assurance/Quality Control Information

Copies of laboratory analytical reports are provided in Appendix C of the attached *2019 Annual Groundwater Monitoring Report*. Laboratory analytical results were reviewed and validated. The data validation and a discussion of any data quality exceptions are provided in Appendix E of the attached *2019 Annual Groundwater Monitoring Report*.

# 10.0 Contour Maps for Each Aquifer Depicting the Potentiometric Gradient for Each Monitoring Event

Groundwater potentiometric surface maps based on the 2019 semiannual gauging results for the shallow saturated zone and the valley fill zone are presented in Figures 4 through 7 of the attached 2019 Annual Groundwater Monitoring Report.

# 11.0 Isoconcentration Maps of Major Constituents of Concern for Each Monitoring Event

The extent of the CGWSL exceedance areas of the following target COCs based on the 2019 semiannual sampling results are presented on Figures 10 through 19 of the attached 2019 Annual Groundwater Monitoring Report: DRO, arsenic, benzene, naphthalene, and MTBE. The extent of the CGWSL exceedance areas of the following water quality parameters based on the 2019 semiannual sampling results are presented on Figures 20 through 29 of the attached 2019 Annual Groundwater Monitoring Report: chloride, fluoride, sulfate, nitrate/nitrite, and TDS.



#### 12.0 PSH Thickness Isopleth Maps for Each Monitoring Event

The presence of PSH and measured PSH thicknesses based on the 2019 semiannual gauging results are shown on Figures 8 and 9 of the attached 2019 Annual Groundwater Monitoring Report.

#### 13.0 Plots of Static Water Elevation Versus Time in Key Wells

Plots presenting PSH thicknesses and static groundwater elevations over time for wells that have historically contained measurable PSH are provided in Appendix D of the attached 2019 Annual Groundwater Monitoring Report.

#### 14.0 Tabulation of the Volumes of PSH Removed

Volumes of groundwater and PSH recovered by the recovery system during 2019 are summarized in Table 5 of the attached 2019 Annual Groundwater Monitoring Report and additional recovery details are provided in Appendix F of the 2019 Annual Groundwater Monitoring Report. An estimated 4,389,902 gallons (104,521 barrels) of groundwater and an estimated 182,749 gallons (4,351 barrels) of PSH were recovered through operation of the automated recovery system in 2019. Further details of the recovery system operation are discussed in Section 6 of the attached 2019 Annual Groundwater Monitoring Report.

#### 15.0 Conclusions and Recommendations

Discharge activities at the refinery during 2019 were conducted in accordance with GW-028. Groundwater conditions at the refinery are generally consistent with previous years. Land application of RO reject water was conducted in accordance with GW-028 from January 1 to January 22, 2019. Land application was discontinued, and after January 22, 2019, RO reject water is further processed in a secondary RO unit, which produces a permeate stream which is utilized in the refinery's cooling towers, and the reject stream from that unit is ultimately discharged to HFNR's injection wells or the City of Artesia's POTW. RO reject water will continue to be utilized in the cooling towers and will not be land applied.



# APPENDIX A Refinery Discharges



# APPENDIX A.1 Refinery Discharges - Treated Wastewater to Injection Wells

# 2019 ANNUAL DISCHARGE PERMIT REPORT HOLLYFRONTIER NAVAJO REFINING LLC - ARTESIA REFINERY DISCHARGE PERMIT GW-028 APPENDIX A.1 SUMMARY OF TREATED WASTEWATER TO INJECTION WELLS

Month	API No. and Well Name	Volume	Average Pressure
(2019)		(bbl)	(psig)
January	30-015-27592 WDW - 1	263,589	1,272
	30-015-20894 WDW - 2	75,463	1,262
	30-015-26575 WDW - 3	92,469	1,177
	30-015-44677 WDW - 4	202,560	58
February	30-015-27592 WDW - 1	188,160	1,270
	30-015-20894 WDW - 2	61,440	1,252
	30-015-26575 WDW - 3	66,240	1,116
	30-015-44677 WDW - 4	228,480	50
March	30-015-27592 WDW - 1	132,857	1,184
	30-015-20894 WDW - 2	53,143	1,157
	30-015-26575 WDW - 3	59,520	1,064
	30-015-44677 WDW - 4	279,531	80
April	30-015-27592 WDW - 1	138,857	1,240
-	30-015-20894 WDW - 2	48,343	1,141
	30-015-26575 WDW - 3	52,457	1,047
	30-015-44677 WDW - 4	240,686	70
May	30-015-27592 WDW - 1	148,800	1,283
	30-015-20894 WDW - 2	62,709	1,215
	30-015-26575 WDW - 3	64,834	1,064
	30-015-44677 WDW - 4	223,200	75
June	30-015-27592 WDW - 1	134,743	1,223
	30-015-20894 WDW - 2	62,743	1,247
	30-015-26575 WDW - 3	55,543	1,031
	30-015-44677 WDW - 4	274,629	98
July	30-015-27592 WDW - 1	139,234	1,219
•	30-015-20894 WDW - 2	60,583	1,224
	30-015-26575 WDW - 3	73,337	1,045
	30-015-44677 WDW - 4	272,091	102
August	30-015-27592 WDW - 1	129,669	1,099
	30-015-20894 WDW - 2	54,206	1,160
	30-015-26575 WDW - 3	81,840	1,062
	30-015-44677 WDW - 4	312,480	115
September	30-015-27592 WDW - 1	126,514	1,133
-	30-015-20894 WDW - 2	48,343	1,137
	30-015-26575 WDW - 3	51,429	993
	30-015-44677 WDW - 4	246,857	101
October	30-015-27592 WDW - 1	125,417	1,117
	30-015-20894 WDW - 2	46,766	1,142
	30-015-26575 WDW - 3	69,086	1,033
	30-015-44677 WDW - 4	247,646	100
November	30-015-27592 WDW - 1	134,743	1,246
	30-015-20894 WDW - 2	45,257	1,227
	30-015-26575 WDW - 3	49,371	1,052
	30-015-44677 WDW - 4	231,429	100
December	30-015-27592 WDW - 1	123,291	1,117
	30-015-20894 WDW - 2	38,263	1,207
	30-015-26575 WDW - 3	53,143	1,097
	30-015-44677 WDW - 4	268,903	109

2019 Cumulative Volume:	bbls
30-015-27592 WDW - 1	1,785,874
30-015-20894 WDW - 2	657,259
30-015-26575 WDW - 3	769,269
30-015-44677 WDW - 4	3,028,492
Total Injected fluids	6,240,894

Average Pressure	psig
30-015-27592 WDW - 1	1,200
30-015-20894 WDW - 2	1,198
30-015-26575 WDW - 3	1,065
30-015-44677 WDW - 4	88

Notes:

API: American Petroleum Institute

bbl: barrel

psig: pounds per square inch gauge



# APPENDIX A.2 Refinery Discharges - Treated Wastewater to City of Artesia POTW

# 2019 ANNUAL DISCHARGE PERMIT REPORT HOLLYFRONTIER NAVAJO REFINING LLC - ARTESIA REFINERY DISCHARGE PERMIT GW-028 APPENDIX A.2

#### SUMMARY OF TREATED WASTEWATER TO THE CITY OF ARTESIA

Refinery WWTP to City of Artesia POTW								
Month (2019)	Rate (gpm)	Volume (gal)						
January	4	178,560						
February	2	89,280						
March	2	89,280						
April	0.1	4,464						
May	0.6	26,784						
June	2.5	111,600						
July	0.4	151,776						
August	3.7	165,168						
September	1.4	62,496						
October	2.3	102,672						
November	2.1	93,744						
December	1.8	80,352						

Average Rate (gpm) 1.91
Cummulative gallons 1,156,176
Cummulative barrels 27,528

#### Notes:

POTW: Publicly-Owned Treatment Works WWTP: Wastewater Treatment Plant

gpm: gallons per minute

gal: gallons

# 2019 ANNUAL DISCHARGE PERMIT REPORT HOLLYFRONTIER NAVAJO REFINING LLC - ARTESIA REFINERY DISCHARGE PERMIT GW-028 APPENDIX A.2

#### SUMMARY OF COOLING TOWER BLOW-DOWN TO THE CITY OF ARTESIA

Cooling Tower Blow-Down to City of Artesia POTW							
Month (2019)	Rate (gpm)	Volume (gal)					
January	73	3,258,720					
February	75	3,348,000					
March	89	3,972,960					
April	92	4,106,880					
May	99	4,419,360					
June	98	4,374,720					
July	104	4,642,560					
August	4,330,080						
September	83	3,705,120					
October	83	3,705,120					
November	63	2,812,320					
December	70	3,124,800					

Average (gpm) 86
Cummulative gallons 45,800,640
Cummulative barrels 1,090,491

Notes:

POTW: Publicly-Owned Treatment Works

gpm: gallons per minute

gal: gallons



# APPENDIX A.3 Refinery Discharges - RO Reject Water Discharge

# 2019 ANNUAL DISCHARGE PERMIT REPORT HOLLYFRONTIER NAVAJO REFINING LLC - ARTESIA REFINERY DISCHARGE PERMIT GW-028 APPENDIX A.3

#### **DAILY AND MONTHLY RO DISCHARGE SUMMARY**

January 2019 - RO Reject Flow/Discharge Measurements								
Skid Location: Measurement:	South Daily Flow	North Daily Flow	Middle Daily Flow	Combined Discharge	Combined Discharge			
Units:	GPM	GPM	GPM	GPM	BPD			
1/1/19	161.66	0.05	162.43	324.13	11,113.03			
1/2/19	162.08	0.04	175.01	337.13	11,558.84			
1/3/19	161.86	0.04	174.48	336.38	11,533.04			
1/4/19	162.07	0.04	174.65	336.76	11,546.17			
1/5/19	161.37	0.05	173.89	335.31	11,496.19			
1/6/19	145.29	0.05	152.94	298.28	10,226.60			
1/7/19	138.63	0.06	141.43	280.13	9,604.41			
1/8/19	162.53	0.06	159.88	322.46	11,055.92			
1/9/19	189.56	0.44	160.14	350.14	12,004.96			
1/10/19	178.74	58.15	95.06	331.95	11,381.00			
1/11/19	170.32	181.47	0.03	351.82	12,062.49			
1/12/19	91.60	180.68	76.75	349.03	11,966.66			
1/13/19	0.00	152.60	175.00	327.60	11,231.94			
1/14/19	0.00	150.41	177.48	327.89	11,241.80			
1/15/19	0.00	151.54	177.72	329.25	11,288.72			
1/16/19	0.00	149.46	174.58	324.05	11,110.12			
1/17/19	0.05	138.39	174.58	313.02	10,732.01			
1/18/19	81.94	76.81	173.34	332.09	11,385.93			
1/19/19	170.69	0.05	173.82	344.56	11,813.42			
1/20/19	141.54	0.05	175.87	317.46	10,884.38			
1/21/19	158.78	0.05	170.51	329.34	11,291.64			
1/22/19*	124.13	0.06	122.70	246.90	8,465.02			
				Total (bbl):	244,994.30			
				Total (gal):	10,289,760.49			

#### Notes:

\* Land application of RO discharge was discontinued after January 22, 2019.

RO: Reverse osmosis

bbl: barrel BPD: barrels per day gal: gallon GPM: gallons per minute



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

February 06, 2019

Scott Denton Navajo Refining Company P.O. Box 159 Artesia, NM 88211-0159

TEL: (575) 748-3311

FAX

RE: RO Reject OrderNo.: 1901787

#### Dear Scott Denton:

Hall Environmental Analysis Laboratory received 2 sample(s) on 1/21/2019 for the analyses presented in the following report.

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

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

ADHS Cert #AZ0682 -- NMED-DWB Cert #NM9425 -- NMED-Micro Cert #NM0901

Sincerely,

Andy Freeman

Laboratory Manager

4901 Hawkins NE

Albuquerque, NM 87109

Date Reported: 2/6/2019

#### Hall Environmental Analysis Laboratory, Inc.

CLIENT: Navajo Refining Company

Client Sample ID: R.O. Reject

 Project:
 RO Reject
 Collection Date: 1/18/2019 1:35:00 PM

 Lab ID:
 1901787-001
 Matrix: AQUEOUS
 Received Date: 1/21/2019 8:20:00 AM

Analyses	Result	PQL	Qual Units	DF	Date Analyzed
EPA METHOD 8011/504.1: EDB					Analyst: <b>JME</b>
1,2-Dibromoethane	ND	0.0093	μg/L	1	1/23/2019 7:52:01 PM
EPA METHOD 8082A: PCB'S					Analyst: <b>TOM</b>
Aroclor 1016	ND	1.0	μg/L	1	1/30/2019 2:45:30 PM
Aroclor 1221	ND	1.0	μg/L	1	1/30/2019 2:45:30 PM
Aroclor 1232	ND	1.0	μg/L	1	1/30/2019 2:45:30 PM
Aroclor 1242	ND	1.0	μg/L	1	1/30/2019 2:45:30 PM
Aroclor 1248	ND	1.0	μg/L	1	1/30/2019 2:45:30 PM
Aroclor 1254	ND	1.0	μg/L	1	1/30/2019 2:45:30 PM
Aroclor 1260	ND	1.0	μg/L	1	1/30/2019 2:45:30 PM
Surr: Decachlorobiphenyl	72.0	24.8-102	%Red	1	1/30/2019 2:45:30 PM
Surr: Tetrachloro-m-xylene	70.4	15.6-106	%Red	1	1/30/2019 2:45:30 PM
EPA METHOD 8015M/D: DIESEL RANGE					Analyst: CLP
Diesel Range Organics (DRO)	ND	1.0	mg/L	1	1/23/2019 9:36:21 AM
Motor Oil Range Organics (MRO)	ND	5.0	mg/L	1	1/23/2019 9:36:21 AM
Surr: DNOP	108	70-130	%Red	1	1/23/2019 9:36:21 AM
EPA METHOD 8310: PAHS					Analyst: <b>TOM</b>
Naphthalene	ND	3.0	μg/L	1	1/30/2019 2:08:37 PM
1-Methylnaphthalene	ND	3.0	μg/L	1	1/30/2019 2:08:37 PM
2-Methylnaphthalene	ND	3.0	μg/L	1	1/30/2019 2:08:37 PM
Acenaphthylene	ND	3.0	μg/L	1	1/30/2019 2:08:37 PM
Acenaphthene	ND	3.0	μg/L	1	1/30/2019 2:08:37 PM
Fluorene	ND	0.80	μg/L	1	1/30/2019 2:08:37 PM
Phenanthrene	ND	0.60	μg/L	1	1/30/2019 2:08:37 PM
Anthracene	ND	0.60	μg/L	1	1/30/2019 2:08:37 PM
Fluoranthene	ND	0.30	μg/L	1	1/30/2019 2:08:37 PM
Pyrene	ND	0.40	μg/L	1	1/30/2019 2:08:37 PM
Benz(a)anthracene	ND	0.070	μg/L	1	1/30/2019 2:08:37 PM
Chrysene	ND	0.20	μg/L	1	1/30/2019 2:08:37 PM
Benzo(b)fluoranthene	ND	0.10	μg/L	1	1/30/2019 2:08:37 PM
Benzo(k)fluoranthene	ND	0.070	μg/L	1	1/30/2019 2:08:37 PM
Benzo(a)pyrene	ND	0.070	μg/L	1	1/30/2019 2:08:37 PM
Dibenz(a,h)anthracene	ND	0.12	μg/L	1	1/30/2019 2:08:37 PM
Benzo(g,h,i)perylene	ND	0.12	μg/L	1	1/30/2019 2:08:37 PM
Indeno(1,2,3-cd)pyrene	ND	0.25	μg/L	1	1/30/2019 2:08:37 PM
Surr: Benzo(e)pyrene	62.8	48.8-93.3	%Red	1	1/30/2019 2:08:37 PM
EPA METHOD 300.0: ANIONS					Analyst: smb
Fluoride	2.3	0.10	mg/L	1	1/21/2019 12:47:38 PM

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

**Qualifiers:** \* Value exceeds Maximum Contaminant Level.

- D Sample Diluted Due to Matrix
- H Holding times for preparation or analysis exceeded
- ND Not Detected at the Reporting Limit
- PQL Practical Quanitative Limit
- S % Recovery outside of range due to dilution or matrix
- B Analyte detected in the associated Method Blank
- E Value above quantitation range
- J Analyte detected below quantitation limits Page 1 of 32
- P Sample pH Not In Range
- RL Reporting Detection Limit
- W Sample container temperature is out of limit as specified

### **Analytical Report**Lab Order **1901787**

#### Hall Environmental Analysis Laboratory, Inc. Date Reported: 2/6/2019

CLIENT: Navajo Refining Company

Client Sample ID: R.O. Reject

 Project:
 RO Reject
 Collection Date: 1/18/2019 1:35:00 PM

 Lab ID:
 1901787-001
 Matrix: AQUEOUS
 Received Date: 1/21/2019 8:20:00 AM

Analyses	Result	PQL Qua	al Units	DF	Date Analyzed
EPA METHOD 300.0: ANIONS					Analyst: smb
Chloride	370	10	mg/L	20	1/21/2019 1:26:13 PM
Sulfate	2000	25	mg/L	50	1/30/2019 12:13:46 AM
Nitrate+Nitrite as N	1.3	1.0	mg/L	5	1/21/2019 6:22:02 PM
EPA METHOD 200.7: DISSOLVED METALS					Analyst: <b>bcv</b>
Aluminum	ND	0.020	mg/L	1	1/24/2019 5:13:02 PM
Barium	0.065	0.0020	mg/L	1	1/24/2019 5:13:02 PM
Beryllium	ND	0.0020	mg/L	1	1/24/2019 5:13:02 PM
Boron	0.11	0.040	mg/L	1	1/24/2019 5:13:02 PM
Cadmium	ND	0.0020	mg/L	1	1/24/2019 5:13:02 PM
Calcium	700	10	mg/L	10	1/24/2019 5:40:37 PM
Chromium	ND	0.0060	mg/L	1	1/24/2019 5:13:02 PM
Cobalt	ND	0.0060	mg/L	1	1/24/2019 5:13:02 PM
Copper	ND	0.0060	mg/L	1	1/24/2019 5:13:02 PM
Iron	ND	0.020	mg/L	1	1/24/2019 5:13:02 PM
Magnesium	220	5.0	mg/L	5	1/24/2019 5:19:53 PM
Manganese	ND	0.0020	mg/L	1	1/24/2019 5:13:02 PM
Molybdenum	ND	0.0080	mg/L	1	1/24/2019 5:13:02 PM
Nickel	ND	0.010	mg/L	1	1/24/2019 5:13:02 PM
Potassium	4.5	1.0	mg/L	1	1/24/2019 5:13:02 PM
Silver	0.010	0.0050	mg/L	1	1/24/2019 5:13:02 PM
Sodium	210	5.0	mg/L	5	1/24/2019 5:19:53 PM
Vanadium	ND	0.050	mg/L	1	1/24/2019 5:13:02 PM
Zinc	0.025	0.010	mg/L	1	1/24/2019 5:13:02 PM
EPA 200.8: DISSOLVED METALS					Analyst: <b>DBK</b>
Antimony	ND	0.0010	mg/L	1	1/24/2019 2:39:19 PM
Arsenic	0.0018	0.0010	mg/L	1	1/24/2019 2:39:19 PM
Lead	ND	0.00050	mg/L	1	1/24/2019 2:39:19 PM
Selenium	0.0090	0.0010	mg/L	1	1/24/2019 2:39:19 PM
Thallium	ND	0.00050	mg/L	1	1/24/2019 2:39:19 PM
Uranium	0.0061	0.00050	mg/L	1	1/24/2019 2:39:19 PM
EPA METHOD 245.1: MERCURY					Analyst: pmf
Mercury	ND	0.00020	mg/L	1	1/24/2019 7:53:16 PM
EPA METHOD 8260B: VOLATILES					Analyst: AG
Benzene	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
Toluene	2.9	1.0	μg/L	1	1/22/2019 5:39:03 PM
Ethylbenzene	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
1,2-Dichloroethane (EDC)	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM

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

**Qualifiers:** \* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

PQL Practical Quanitative Limit

S % Recovery outside of range due to dilution or matrix

- B Analyte detected in the associated Method Blank
- E Value above quantitation range
- J Analyte detected below quantitation limits Page 2 of 32
- P Sample pH Not In Range
- RL Reporting Detection Limit
- W Sample container temperature is out of limit as specified

#### **Analytical Report**

### Lab Order **1901787**Date Reported: **2/6/2019**

#### Hall Environmental Analysis Laboratory, Inc.

CLIENT: Navajo Refining Company

Client Sample ID: R.O. Reject

 Project:
 RO Reject
 Collection Date: 1/18/2019 1:35:00 PM

 Lab ID:
 1901787-001
 Matrix: AQUEOUS
 Received Date: 1/21/2019 8:20:00 AM

Analyses	Result	PQL Qua	al Units	DF	Date Analyzed
EPA METHOD 8260B: VOLATILES					Analyst: AG
1,2-Dibromoethane (EDB)	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
Carbon Tetrachloride	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
Chloroform	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
1,1-Dichloroethane	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
1,1-Dichloroethene	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
Methylene Chloride	ND	3.0	μg/L	1	1/22/2019 5:39:03 PM
1,1,2,2-Tetrachloroethane	ND	2.0	μg/L	1	1/22/2019 5:39:03 PM
Tetrachloroethene (PCE)	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
1,1,1-Trichloroethane	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
1,1,2-Trichloroethane	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
Trichloroethene (TCE)	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
Vinyl chloride	ND	1.0	μg/L	1	1/22/2019 5:39:03 PM
Xylenes, Total	ND	1.5	μg/L	1	1/22/2019 5:39:03 PM
Surr: 1,2-Dichloroethane-d4	110	70-130	%Rec	1	1/22/2019 5:39:03 PM
Surr: 4-Bromofluorobenzene	107	70-130	%Rec	1	1/22/2019 5:39:03 PM
Surr: Dibromofluoromethane	108	70-130	%Rec	1	1/22/2019 5:39:03 PM
Surr: Toluene-d8	104	70-130	%Rec	1	1/22/2019 5:39:03 PM
EPA METHOD 8015D: GASOLINE RANGE					Analyst: AG
Gasoline Range Organics (GRO)	ND	0.050	mg/L	1	1/22/2019 5:39:03 PM
Surr: BFB	99.7	70-130	%Rec	1	1/22/2019 5:39:03 PM
TOTAL PHENOLICS BY SW-846 9067					Analyst: CLP
Phenolics	ND	2.5	μg/L	1	1/30/2019
EPA 8270D: SEMIVOLATILES					Analyst: PAC
Atrazine	ND	0.010	μg/L	1	1/25/2019
1,2,4-Trichlorobenzene	ND	0.010	μg/L	1	1/25/2019
2,4,6-Trichlorophenol	ND	0.010	μg/L	1	1/25/2019
2,4-Dichlorophenol	ND	0.010	μg/L	1	1/25/2019
2,4-Dimethylphenol	ND	0.010	μg/L	1	1/25/2019
2,4-Dinitrophenol	ND	0.010	μg/L	1	1/25/2019
2,4-Dinitrotoluene	ND	0.010	μg/L	1	1/25/2019
2,6-Dinitrotoluene	ND	0.010	μg/L	1	1/25/2019
2-Chloronaphthalene	ND	0.0010	μg/L	1	1/25/2019
2-Chlorophenol	ND	0.010	μg/L	1	1/25/2019
2-Nitrophenol	ND	0.010	μg/L	1	1/25/2019
3,3´-Dichlorobenzidine	ND	0.010	μg/L	1	1/25/2019
4,6-Dinitro-2-methylphenol	ND	0.010	μg/L	1	1/25/2019
4-Bromophenyl phenyl ether	ND	0.010	μg/L	1	1/25/2019
4-Chloro-3-methylphenol	ND	0.010	μg/L	1	1/25/2019

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

- \* Value exceeds Maximum Contaminant Level.
- D Sample Diluted Due to Matrix
- H Holding times for preparation or analysis exceeded
- ND Not Detected at the Reporting Limit
- PQL Practical Quanitative Limit
- S % Recovery outside of range due to dilution or matrix
- B Analyte detected in the associated Method Blank
- E Value above quantitation range
- J Analyte detected below quantitation limits Page 3 of 32
- P Sample pH Not In Range
- RL Reporting Detection Limit
- W Sample container temperature is out of limit as specified

#### **Analytical Report**

### Lab Order **1901787**Date Reported: **2/6/2019**

#### Hall Environmental Analysis Laboratory, Inc.

CLIENT: Navajo Refining Company

Client Sample ID: R.O. Reject

 Project:
 RO Reject
 Collection Date: 1/18/2019 1:35:00 PM

 Lab ID:
 1901787-001
 Matrix: AQUEOUS
 Received Date: 1/21/2019 8:20:00 AM

Analyses	Result	PQL Qua	al Units	DF	Date Analyzed
EPA 8270D: SEMIVOLATILES					Analyst: PAC
4-Chlorophenyl phenyl ether	ND	0.010	μg/L	1	1/25/2019
4-Nitrophenol	ND	0.010	μg/L	1	1/25/2019
Acenaphthene	ND	0.0010	μg/L	1	1/25/2019
Acenaphthylene	ND	0.0010	μg/L	1	1/25/2019
Anthracene	ND	0.0010	μg/L	1	1/25/2019
Benzidine	ND	0.010	μg/L	1	1/25/2019
Benzo(g,h,i)perylene	ND	0.0010	μg/L	1	1/25/2019
Benz(a)anthracene	ND	0.0010	μg/L	1	1/25/2019
Benzo(a)pyrene	ND	0.0010	μg/L	1	1/25/2019
Benzo(b)fluoranthene	ND	0.0010	μg/L	1	1/25/2019
Benzo(k)fluoranthene	ND	0.0010	μg/L	1	1/25/2019
Bis(2-chloroethoxy)methane	ND	0.010	μg/L	1	1/25/2019
Bis(2-chloroethyl)ether	ND	0.010	μg/L	1	1/25/2019
Bis(2-chloroisopropyl)ether	ND	0.010	μg/L	1	1/25/2019
Bis(2-ethylhexyl)phthalate	ND	0.0030	μg/L	1	1/25/2019
Butyl benzyl phthalate	ND	0.0030	μg/L	1	1/25/2019
Chrysene	ND	0.0010	μg/L	1	1/25/2019
Dibenz(a,h)anthracene	ND	0.0010	μg/L	1	1/25/2019
Diethyl phthalate	ND	0.0030	μg/L	1	1/25/2019
Dimethyl phthalate	ND	0.0030	μg/L	1	1/25/2019
Di-n-butyl phthalate	ND	0.0030	μg/L	1	1/25/2019
Di-n-octyl phthalate	ND	0.0030	μg/L	1	1/25/2019
Fluoranthene	ND	0.0010	μg/L	1	1/25/2019
Fluorene	ND	0.0010	μg/L	1	1/25/2019
Hexachlorobenzene	ND	0.0010	μg/L	1	1/25/2019
Hexachlorobutadiene	ND	0.010	μg/L	1	1/25/2019
Hexachlorocyclopentadiene	ND	0.010	μg/L	1	1/25/2019
Hexachloroethane	ND	0.010	μg/L	1	1/25/2019
Indeno(1,2,3-cd)pyrene	ND	0.0010	μg/L	1	1/25/2019
Isophorone	ND	0.010	μg/L	1	1/25/2019
Naphthalene	ND	0.0010	μg/L	1	1/25/2019
Nitrobenzene	ND	0.010	μg/L	1	1/25/2019
N-Nitrosodimethylamine	ND	0.010	μg/L	1	1/25/2019
N-Nitrosodi-n-propylamine	ND	0.010	μg/L	1	1/25/2019
N-Nitrosodiphenylamine	ND	0.010	μg/L	1	1/25/2019
Pentachlorophenol	ND	0.010	μg/L	1	1/25/2019
Phenanthrene	ND	0.0010	μg/L	1	1/25/2019
Phenol	ND	0.010	μg/L	1	1/25/2019
Pyrene	ND	0.0010	μg/L	1	1/25/2019

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

- \* Value exceeds Maximum Contaminant Level.
- D Sample Diluted Due to Matrix
- H Holding times for preparation or analysis exceeded
- ND Not Detected at the Reporting Limit
- PQL Practical Quanitative Limit
- S % Recovery outside of range due to dilution or matrix
- B Analyte detected in the associated Method Blank
- E Value above quantitation range
- J Analyte detected below quantitation limits Page 4 of 32
- P Sample pH Not In Range
- RL Reporting Detection Limit
- W Sample container temperature is out of limit as specified

### **Analytical Report**Lab Order **1901787**

#### Hall Environmental Analysis Laboratory, Inc.

Date Reported: 2/6/2019

CLIENT: Navajo Refining Company

Client Sample ID: R.O. Reject

 Project:
 RO Reject
 Collection Date: 1/18/2019 1:35:00 PM

 Lab ID:
 1901787-001
 Matrix: AQUEOUS
 Received Date: 1/21/2019 8:20:00 AM

Analyses	Result	PQL Qu	ual Units	DF	Date Analyzed
EPA 335.4: TOTAL CYANIDE SUBBED					Analyst: PAC
Cyanide	ND	0.00500	mg/L	1	1/29/2019
EPA 903.1: RA 226 AND EPA 904.0: RA 228-SUE	BED				Analyst: PAC
Radium-226	2.21	0.741	pCi/L	1	1/29/2019
Radium-226 ±	0.903	0.741	pCi/L	1	1/29/2019
Radium-228	0.0923	0.645	pCi/L	1	1/29/2019
Radium-228 ±	0.288	0.645	pCi/L	1	1/29/2019
SM2510B: SPECIFIC CONDUCTANCE					Analyst: MRA
Conductivity	4300	5.0	µmhos/	c 1	1/21/2019 4:11:55 PM
SM4500-H+B / 9040C: PH					Analyst: MRA
рН	8.02		H pH units	s 1	1/21/2019 4:11:55 PM
SM2540C MOD: TOTAL DISSOLVED SOLIDS					Analyst: KS
Total Dissolved Solids	4020	20.0	* mg/L	1	1/23/2019 3:44:00 PM

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

- Value exceeds Maximum Contaminant Level.
- D Sample Diluted Due to Matrix
- H Holding times for preparation or analysis exceeded
- ND Not Detected at the Reporting Limit
- PQL Practical Quanitative Limit
- S % Recovery outside of range due to dilution or matrix
- B Analyte detected in the associated Method Blank
- E Value above quantitation range
- J Analyte detected below quantitation limits Page 5 of 32
- P Sample pH Not In Range
- RL Reporting Detection Limit
- W Sample container temperature is out of limit as specified

### **Analytical Report**Lab Order **1901787**

#### Hall Environmental Analysis Laboratory, Inc.

Date Reported: 2/6/2019

CLIENT: Navajo Refining Company Client Sample ID: Trip Blank

**Project:** RO Reject Collection Date:

**Lab ID:** 1901787-002 **Matrix:** TRIP BLANK **Received Date:** 1/21/2019 8:20:00 AM

Analyses	Result	PQL Qu	al Units	DF	Date Analyzed
EPA METHOD 8011/504.1: EDB					Analyst: <b>JME</b>
1,2-Dibromoethane	ND	0.0097	μg/L	1	1/23/2019 8:51:15 PM
EPA METHOD 8260B: VOLATILES					Analyst: DJF
Benzene	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
Toluene	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
Ethylbenzene	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
1,2-Dichloroethane (EDC)	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
1,2-Dibromoethane (EDB)	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
Carbon Tetrachloride	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
Chloroform	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
1,1-Dichloroethane	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
1,1-Dichloroethene	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
Methylene Chloride	ND	3.0	μg/L	1	1/23/2019 6:30:47 AM
1,1,2,2-Tetrachloroethane	ND	2.0	μg/L	1	1/23/2019 6:30:47 AM
Tetrachloroethene (PCE)	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
1,1,1-Trichloroethane	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
1,1,2-Trichloroethane	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
Trichloroethene (TCE)	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
Vinyl chloride	ND	1.0	μg/L	1	1/23/2019 6:30:47 AM
Xylenes, Total	ND	1.5	μg/L	1	1/23/2019 6:30:47 AM
Surr: 1,2-Dichloroethane-d4	103	70-130	%Rec	1	1/23/2019 6:30:47 AM
Surr: 4-Bromofluorobenzene	101	70-130	%Rec	1	1/23/2019 6:30:47 AM
Surr: Dibromofluoromethane	138	70-130	S %Rec	1	1/23/2019 6:30:47 AM
Surr: Toluene-d8	103	70-130	%Rec	1	1/23/2019 6:30:47 AM

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

- \* Value exceeds Maximum Contaminant Level.
- D Sample Diluted Due to Matrix
- H Holding times for preparation or analysis exceeded
- ND Not Detected at the Reporting Limit
- PQL Practical Quanitative Limit
- S % Recovery outside of range due to dilution or matrix
- B Analyte detected in the associated Method Blank
- E Value above quantitation range
- J Analyte detected below quantitation limits Page 6 of 32
- P Sample pH Not In Range
- RL Reporting Detection Limit
- W Sample container temperature is out of limit as specified

#### Hall Environmental Analysis Laboratory, Inc.

1.0

0.050

0.010

ND

ND

WO#: 1901787

06-Feb-19

**Client:** Navajo Refining Company

**Project: RO** Reject

Sample ID MB-B SampType: MBLK TestCode: EPA Method 200.7: Dissolved Metals Client ID: PBW Batch ID: **B57243** RunNo: 57243 Prep Date: Analysis Date: 1/24/2019 SeqNo: 1914910 Units: mg/L Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Aluminum ND 0.020 0.0020 ND Barium Beryllium ND 0.0020 Boron ND 0.040 Cadmium ND 0.0020 Calcium ND 1.0 Chromium ND 0.0060 ND 0.0060 Cobalt Copper ND 0.0060 ND 0.020 Iron Magnesium ND 1.0 ND 0.0020 Manganese 0.0080 Molybdenum ND Nickel ND 0.010 Potassium ND 1.0 Silver ND 0.0050 Sodium ND

Sample ID LLLCS-B	SampType: LCSLL TestCode: EPA M					PA Method	200.7: Dissol	ved Metal	ls	
Client ID: BatchQC	Bato	ch ID: B5	7243	R	RunNo: 57243					
Prep Date:	Analysis	Date: 1/	24/2019	S	SeqNo: 1	914911	Units: mg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Aluminum	ND	0.020	0.01000	0	129	50	150			
Barium	ND	0.0020	0.002000	0	94.4	50	150			
Beryllium	ND	0.0020	0.002000	0	99.7	50	150			
Boron	ND	0.040	0.04000	0	97.1	50	150			
Cadmium	ND	0.0020	0.002000	0	67.2	50	150			
Calcium	ND	1.0	0.5000	0	104	50	150			
Chromium	ND	0.0060	0.006000	0	92.2	50	150			
Cobalt	0.0065	0.0060	0.006000	0	108	50	150			
Copper	ND	0.0060	0.006000	0	68.8	50	150			
Iron	0.021	0.020	0.02000	0	105	50	150			
Magnesium	ND	1.0	0.5000	0	103	50	150			
Manganese	ND	0.0020	0.002000	0	96.5	50	150			
Molybdenum	ND	0.0080	0.008000	0	80.0	50	150			
Nickel	ND	0.010	0.005000	0	70.3	50	150			

#### Qualifiers:

Vanadium

Zinc

- Value exceeds Maximum Contaminant Level.
- D Sample Diluted Due to Matrix
- Η Holding times for preparation or analysis exceeded
- ND Not Detected at the Reporting Limit
- POL Practical Quanitative Limit
- % Recovery outside of range due to dilution or matrix
- В Analyte detected in the associated Method Blank
- Е Value above quantitation range

Reporting Detection Limit

- J Analyte detected below quantitation limits
- P Sample pH Not In Range

RL

Sample container temperature is out of limit as specified

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#### Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID LLLCS-B	SampType: LCSLL			TestCode: EPA Method 200.7: Dissolved Metals						
Client ID: BatchQC	Batch ID: <b>B57243</b>			F	RunNo: 5					
Prep Date:	Analysis	Date: 1/	24/2019	/2019 SeqNo: 1914			914911 Units: mg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Potassium	ND	1.0	0.5000	0	104	50	150			
Silver	ND	0.0050	0.005000	0	83.4	50	150			
Sodium	ND	1.0	0.5000	0	117	50	150			
Vanadium	ND	0.050	0.01000	0	88.0	50	150			
Zinc	ND	0.010	0.005000	0	126	50	150			

Sample ID LCS-B SampType: LCS			S	TestCode: EPA Method 200.7: Dissolved Metals								
Client ID: LCSW	Batch ID: <b>B57243</b> Analysis Date: <b>1/24/2019</b>			RunNo: <b>57243</b>								
Prep Date:				SeqNo: 1914912			Units: mg/L					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual		
Aluminum	0.55	0.020	0.5000	0	109	85	115					
Barium	0.49	0.0020	0.5000	0	98.2	85	115					
Beryllium	0.50	0.0020	0.5000	0	99.3	85	115					
Boron	0.51	0.040	0.5000	0	102	85	115					
Cadmium	0.50	0.0020	0.5000	0	99.6	85	115					
Calcium	50	1.0	50.00	0	99.0	85	115					
Chromium	0.49	0.0060	0.5000	0	97.8	85	115					
Cobalt	0.48	0.0060	0.5000	0	95.5	85	115					
Copper	0.50	0.0060	0.5000	0	99.3	85	115					
Iron	0.49	0.020	0.5000	0	97.6	85	115					
Magnesium	50	1.0	50.00	0	100	85	115					
Manganese	0.48	0.0020	0.5000	0	96.8	85	115					
Molybdenum	0.49	0.0080	0.5000	0	98.5	85	115					
Nickel	0.49	0.010	0.5000	0	97.1	85	115					
Potassium	50	1.0	50.00	0	99.3	85	115					
Silver	0.10	0.0050	0.1000	0	102	85	115					
Sodium	51	1.0	50.00	0	101	85	115					
Vanadium	0.50	0.050	0.5000	0	99.1	85	115					
Zinc	0.48	0.010	0.5000	0	96.2	85	115					

Sample ID 1901787-001GMS	Samp	SampType: <b>MS</b>				TestCode: EPA Method 200.7: Dissolved Metals						
Client ID: R.O. Reject	Bato	h ID: <b>B5</b>	7243	F	RunNo: 5							
Prep Date:	Analysis I	Date: 1/	24/2019	8	SeqNo: 1	915074	Units: mg/L					
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual		
Aluminum	0.55	0.020	0.5000	0	109	70	130					
Barium	0.52	0.0020	0.5000	0.06527	91.7	70	130					
Beryllium	0.52	0.0020	0.5000	0.0003350	103	70	130					
Boron	0.61	0.040	0.5000	0.1099	100	70	130					

#### Qualifiers:

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

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#### Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID 1901787-001GMS SampType: MS TestCode: EPA Method 200.7: Dissolved Metals Client ID: R.O. Reject Batch ID: **B57243** RunNo: 57243 Prep Date: Analysis Date: 1/24/2019 SeqNo: 1915074 Units: mg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Cadmium 0.48 0.0020 0.5000 95.6 0 70 130 87.2 70 Chromium 0.44 0.0060 0.5000 0 130 89.7 Cobalt 0.45 0.0060 0.5000 0 70 130 Copper 0.52 0.0060 0.5000 0.002356 103 70 130 Iron 0.50 0.020 0.5000 0 100 70 130 0.48 0.0020 0.5000 0 96.3 70 130 Manganese Molybdenum 0.44 0.0080 0.5000 0 88.3 70 130 0.45 0.010 0.5000 0 89.9 70 Nickel 130 Potassium 53 1.0 50.00 4.528 96.9 70 130 0.0050 80.3 Silver 0.091 0.1000 0.01047 70 130 Vanadium 0.48 0.050 0.5000 0.01386 93.7 70 130 0.47 0.010 0.5000 0.02475 89.1 70 130 Zinc

Sample ID 1901787-001GMS	TestCode: EPA Method 200.7: Dissolved Metals									
Client ID: R.O. Reject	Batch ID: <b>B57243</b>			RunNo: <b>57243</b>						
Prep Date:	Analysis	Date: 1/	24/2019	9	SeqNo: 1	915075	Units: mg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Aluminum	0.55	0.020	0.5000	0	110	70	130	0.461	20	•
Barium	0.53	0.0020	0.5000	0.06527	93.0	70	130	1.20	20	
Beryllium	0.52	0.0020	0.5000	0.0003350	103	70	130	0.258	20	
Boron	0.62	0.040	0.5000	0.1099	102	70	130	1.46	20	
Cadmium	0.48	0.0020	0.5000	0	97.0	70	130	1.40	20	
Chromium	0.44	0.0060	0.5000	0	87.9	70	130	0.767	20	
Cobalt	0.45	0.0060	0.5000	0	90.5	70	130	0.924	20	
Copper	0.53	0.0060	0.5000	0.002356	105	70	130	1.42	20	
Iron	0.51	0.020	0.5000	0	103	70	130	2.28	20	
Manganese	0.49	0.0020	0.5000	0	97.1	70	130	0.820	20	
Molybdenum	0.45	0.0080	0.5000	0	89.5	70	130	1.30	20	
Nickel	0.46	0.010	0.5000	0	91.2	70	130	1.48	20	
Potassium	55	1.0	50.00	4.528	101	70	130	4.22	20	
Silver	0.092	0.0050	0.1000	0.01047	81.0	70	130	0.802	20	
Vanadium	0.49	0.050	0.5000	0.01386	94.7	70	130	1.01	20	
Zinc	0.47	0.010	0.5000	0.02475	90.0	70	130	0.963	20	

Sample ID 1901787-001GMS SampType: MS TestCode: EPA Method 200.7: Dissolved Metals

Client ID: R.O. Reject Batch ID: B57243 RunNo: 57243

Prep Date: Analysis Date: 1/24/2019 SeqNo: 1915077 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

#### Qualifiers:

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

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#### Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

**Client:** Navajo Refining Company

**Project:** RO Reject

Sample ID 1901787-001GMS SampType: MS TestCode: EPA Method 200.7: Dissolved Metals

Client ID: R.O. Reject Batch ID: B57243 RunNo: 57243

Prep Date: Analysis Date: 1/24/2019 SeqNo: 1915077 Units: mg/L

Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Magnesium 470 250.0 216.0 100 70 5.0 130 Sodium 470 250.0 70 5.0 212.8 103 130

Sample ID 1901787-001GMSD SampType: MSD TestCode: EPA Method 200.7: Dissolved Metals

Client ID: R.O. Reject Batch ID: B57243 RunNo: 57243

Prep Date: Analysis Date: 1/24/2019 SeqNo: 1915078 Units: mg/L

LowLimit Analyte Result **PQL** SPK value SPK Ref Val %REC HighLimit %RPD **RPDLimit** Qual Magnesium 470 5.0 250.0 216.0 102 70 130 0.864 20 480 250.0 107 70 20 Sodium 5.0 212.8 130 1.85

#### Qualifiers:

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

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## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

**Client:** Navajo Refining Company

**Project:** RO Reject

Sample ID MB SampType: MBLK TestCode: EPA 200.8: Dissolved Metals Client ID: PBW Batch ID: **B57230** RunNo: 57230 Prep Date: Analysis Date: 1/24/2019 SeqNo: 1914365 Units: mg/L Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual ND 0.0010 Antimony 0.0010 Arsenic ND 0.00050 Lead ND Selenium ND 0.0010 Thallium ND 0.00050 Uranium ND 0.00050

Sample ID MSLLLCS	Samp	Type: <b>LC</b>	SLL	Tes	tCode: E	PA 200.8: I	Dissolved Me	tals		
Client ID: BatchQC	Bate	ch ID: <b>B5</b>	7230	F	RunNo: <b>5</b>	7230				
Prep Date:	Analysis	Date: 1/	24/2019	S	SeqNo: 1	914366	Units: mg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Antimony	ND	0.0010	0.001000	0	94.3	50	150			
Arsenic	0.0011	0.0010	0.001000	0	108	50	150			
Lead	ND	0.00050	0.0005000	0	94.7	50	150			
Selenium	0.0011	0.0010	0.001000	0	106	50	150			
Thallium	ND	0.00050	0.0005000	0	93.4	50	150			
Uranium	ND	0.00050	0.0005000	0	94.2	50	150			

Sample ID MSLCS	SampType: LC	S	Tes	tCode: El	PA 200.8: [	Dissolved Met	als		
Client ID: LCSW	Batch ID: B5	7230	F	RunNo: 5	7230				
Prep Date:	Analysis Date: 1/	/24/2019	8	SeqNo: 1	914367	Units: mg/L			
Analyte	Result PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Antimony	0.023 0.0010	0.02500	0	92.7	85	115			
Arsenic	0.024 0.0010	0.02500	0	95.4	85	115			
Lead	0.012 0.00050	0.01250	0	94.0	85	115			
Selenium	0.024 0.0010	0.02500	0	95.8	85	115			
Thallium	0.012 0.00050	0.01250	0	94.3	85	115			
Uranium	0.012 0.00050	0.01250	0	93.9	85	115			

#### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

PQL Practical Quanitative Limit

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

Sample pH Not In Range

RL Reporting Detection Limit

P

W Sample container temperature is out of limit as specified

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#### Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

**Client:** Navajo Refining Company

**Project:** RO Reject

Sample ID MB-42793 SampType: MBLK TestCode: EPA Method 245.1: Mercury

Client ID: PBW Batch ID: 42793 RunNo: 57245

Prep Date: 1/24/2019 Analysis Date: 1/24/2019 SeqNo: 1914854 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Mercury ND 0.00020

Sample ID LCS-42793 SampType: LCS TestCode: EPA Method 245.1: Mercury

Client ID: LCSW Batch ID: 42793 RunNo: 57245

Prep Date: 1/24/2019 Analysis Date: 1/24/2019 SeqNo: 1914855 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Mercury 0.0049 0.00020 0.005000 0 98.9 80 120

Sample ID 1901832-001CMS SampType: MS TestCode: EPA Method 245.1: Mercury

Client ID: BatchQC Batch ID: 42793 RunNo: 57245

Prep Date: 1/24/2019 Analysis Date: 1/24/2019 SeqNo: 1914859 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Mercury 0.0049 0.00020 0.005000 0 98.3 75 125

Sample ID 1901832-001CMSD SampType: MSD TestCode: EPA Method 245.1: Mercury

Client ID: BatchQC Batch ID: 42793 RunNo: 57245

Prep Date: 1/24/2019 Analysis Date: 1/24/2019 SeqNo: 1914860 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Mercury 0.0048 0.00020 0.005000 0 95.2 75 125 3.26 20

#### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

PQL Practical Quanitative Limit

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

D G 1 HN I D

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P Sample pH Not In Range

RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

#### Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID MB SampType: MBLK TestCode: EPA Method 300.0: Anions

Client ID: PBW Batch ID: R57149 RunNo: 57149

Olicit ID. 1 Day Daton ID. N3/143 Null No. 3/143

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1911765 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

 Fluoride
 ND
 0.10

 Chloride
 ND
 0.50

 Nitrate+Nitrite as N
 ND
 0.20

Sample ID LCS SampType: LCS TestCode: EPA Method 300.0: Anions

Client ID: LCSW Batch ID: R57149 RunNo: 57149

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1911766 Units: mg/L

Analyte **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Fluoride 0.48 0.10 O 96.4 90 110 0.5000 Chloride 4.8 0.50 5.000 0 95.5 90 110 0.20 0 99.1 90 Nitrate+Nitrite as N 3.5 3.500 110

Sample ID 1901787-001EMS SampType: MS TestCode: EPA Method 300.0: Anions

Client ID: R.O. Reject Batch ID: R57149 RunNo: 57149

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1911772 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Fluoride 3.0 0.10 0.5000 2.347 125 66.7 127

Sample ID 1901787-001EMSD SampType: MSD TestCode: EPA Method 300.0: Anions

Client ID: R.O. Reject Batch ID: R57149 RunNo: 57149

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1911773 Units: mg/L

**RPDLimit** SPK value SPK Ref Val Analyte Result POL %REC LowLimit HighLimit %RPD Qual Fluoride 3.0 0.10 0.5000 2.347 125 66.7 127 0.0950 20

Sample ID 1901792-001BMS SampType: MS TestCode: EPA Method 300.0: Anions Client ID: **BatchQC** Batch ID: R57149 RunNo: 57149 Prep Date: Analysis Date: 1/21/2019 SeqNo: 1911802 Units: mg/L SPK value SPK Ref Val %REC %RPD **RPDLimit** Result **PQL** LowLimit HighLimit Qual

Analyte Fluoride 0.83 0.10 0.5000 0.3410 98.4 66.7 127 Nitrate+Nitrite as N 5.6 0.20 3.500 2.098 100 70 117

Sample ID 1901792-001BMSD SampType: MSD TestCode: EPA Method 300.0: Anions

Client ID: BatchQC Batch ID: R57149 RunNo: 57149

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1911803 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

#### Qualifiers:

\* Value exceeds Maximum Contaminant Level. B Analyte detected in the associated Method Blank

D Sample Diluted Due to Matrix E Value above quantitation range

H Holding times for preparation or analysis exceeded J Analyte detected below quantitation limits Page 13 of 32

ND Not Detected at the Reporting Limit P Sample pH Not In Range
PQL Practical Quanitative Limit RL Reporting Detection Limit

S % Recovery outside of range due to dilution or matrix W Sample container temperature is out of limit as specified

#### Hall Environmental Analysis Laboratory, Inc.

WO#: 1901787

06-Feb-19

**Client:** Navajo Refining Company

**Project: RO** Reject

Sample ID 1901792-001BMSD SampType: MSD TestCode: EPA Method 300.0: Anions

Client ID: **BatchQC** Batch ID: **R57149** RunNo: 57149

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1911803 Units: mg/L

Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Fluoride 0.10 98.3 66.7 0.83 0.5000 0.3410 127 0.0414 20 Nitrate+Nitrite as N 5.6 0.20 3.500 2.098 101 70 117 0.447 20

Sample ID MB SampType: MBLK TestCode: EPA Method 300.0: Anions Client ID: PBW Batch ID: **R57344** RunNo: 57344 Prep Date: Analysis Date: 1/29/2019 SeqNo: 1918656 Units: mg/L %REC LowLimit Analyte Result POL SPK value SPK Ref Val HighLimit %RPD **RPDLimit** Qual

Sulfate ND 0.50

Sample ID LCS SampType: LCS TestCode: EPA Method 300.0: Anions

Client ID: LCSW Batch ID: **R57344** RunNo: 57344

Prep Date: Analysis Date: 1/29/2019 SeqNo: 1918657 Units: mg/L

Analyte **PQL** SPK value SPK Ref Val %REC HighLimit %RPD **RPDLimit** Qual LowLimit

0.50 90 Sulfate 10 10.00 0 102 110

Sample ID 1901A59-001BMS SampType: MS TestCode: EPA Method 300.0: Anions

Client ID: **BatchQC** Batch ID: **R57344** RunNo: 57344

Prep Date: Analysis Date: 1/29/2019 SeqNo: 1918659 Units: mg/L

**PQL** %REC %RPD **RPDLimit** Analyte Result SPK value SPK Ref Val LowLimit HighLimit Qual

Sulfate 45 0.50 10.00 33.48 117 74.9 123

Sample ID 1901A59-001BMSD SampType: MSD TestCode: EPA Method 300.0: Anions

Client ID: **BatchQC** Batch ID: **R57344** RunNo: 57344

Prep Date: Analysis Date: 1/29/2019 SeqNo: 1918660 Units: mq/L

Analyte SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Result PQI Qual 33.48 Sulfate 45 0.50 10.00 120 74.9 123 0.623

#### Qualifiers:

Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

Η Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

POL Practical Quanitative Limit

% Recovery outside of range due to dilution or matrix

В Analyte detected in the associated Method Blank

Е Value above quantitation range

J Analyte detected below quantitation limits

Page 14 of 32

P Sample pH Not In Range

RLReporting Detection Limit

Sample container temperature is out of limit as specified

## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

Qual

%RPD

HighLimit

**RPDLimit** 

06-Feb-19

**Client:** Navajo Refining Company

**Project:** RO Reject

Analyte

Sample ID MB-42735 SampType: MBLK TestCode: EPA Method 8011/504.1: EDB

Client ID: PBW Batch ID: 42735 RunNo: 57205

Prep Date: 1/23/2019 Analysis Date: 1/23/2019 SeqNo: 1913435 Units: μg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

SPK value SPK Ref Val %REC LowLimit

1,2-Dibromoethane ND 0.010

Sample ID LCS-42735 SampType: LCS TestCode: EPA Method 8011/504.1: EDB

Client ID: LCSW Batch ID: 42735 RunNo: 57205

PQL

Result

Prep Date: 1/23/2019 Analysis Date: 1/23/2019 SeqNo: 1913437 Units: μg/L

Trep Date. 1723/2019 Analysis Date. 1723/2019 Sequit. 1913437 Offics. Hg/L

1,2-Dibromoethane 0.085 0.010 0.1000 0 85.3 70 130

Sample ID 1901787-001BMS SampType: MS TestCode: EPA Method 8011/504.1: EDB

Client ID: R.O. Reject Batch ID: 42735 RunNo: 57205

Prep Date: 1/23/2019 Analysis Date: 1/23/2019 SeqNo: 1913497 Units: μg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

1.2-Dibromoethane 0.067 0.0093 0.09333 0 72.2 55 129

Sample ID 1901787-001BMSD SampType: MSD TestCode: EPA Method 8011/504.1: EDB

Client ID: R.O. Reject Batch ID: 42735 RunNo: 57205

Prep Date: 1/23/2019 Analysis Date: 1/23/2019 SeqNo: 1913499 Units: μg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

1,2-Dibromoethane 0.061 0.0093 0.09333 0 65.0 55 125 10.4 20

#### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

PQL Practical Quanitative Limit

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

P Sample pH Not In Range

RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

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## Hall Environmental Analysis Laboratory, Inc.

Result

5.7

0.50

**PQL** 

1.0

WO#: 1901787

06-Feb-19

**Client:** Navajo Refining Company

**Project: RO** Reject

Sample ID MB-42745 SampType: MBLK TestCode: EPA Method 8015M/D: Diesel Range Client ID: PBW Batch ID: 42745 RunNo: 57173 Prep Date: 1/22/2019 Analysis Date: 1/23/2019 SeqNo: 1913176 Units: mg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Diesel Range Organics (DRO) ND 1.0 Motor Oil Range Organics (MRO) ND 5.0 Surr: DNOP 70 0.98 1.000 98.2 130 Sample ID LCS-42745 SampType: LCS TestCode: EPA Method 8015M/D: Diesel Range Client ID: LCSW Batch ID: 42745 RunNo: 57173 Prep Date: Analysis Date: 1/23/2019 SeqNo: 1913177 1/22/2019 Units: mg/L Analyte **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Diesel Range Organics (DRO) 5.6 1.0 5.000 112 71.8 135 Surr: DNOP 0.50 0.5000 99.8 70 130 Sample ID 1901789-001BMS SampType: MS TestCode: EPA Method 8015M/D: Diesel Range Client ID: **BatchQC** Batch ID: 42745 RunNo: 57173 Prep Date: 1/22/2019 Analysis Date: 1/23/2019 SeqNo: 1913184 Units: mg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Diesel Range Organics (DRO) 5.6 0 112 68.1 1.0 5.000 137 Surr: DNOP 0.50 0.5000 99.3 70 130 Sample ID 1901789-001BMSD SampType: MSD TestCode: EPA Method 8015M/D: Diesel Range Client ID: **BatchQC** Batch ID: 42745 RunNo: 57173 Prep Date: 1/22/2019 Analysis Date: 1/23/2019 SeqNo: 1913185 Units: mg/L

SPK value SPK Ref Val

0

5.000

0.5000

#### Qualifiers:

Analyte

Surr: DNOP

Diesel Range Organics (DRO)

- Value exceeds Maximum Contaminant Level.
- D Sample Diluted Due to Matrix
- Η Holding times for preparation or analysis exceeded
- ND Not Detected at the Reporting Limit
- POL Practical Quanitative Limit
- % Recovery outside of range due to dilution or matrix
- В Analyte detected in the associated Method Blank

LowLimit

68.1

70

HighLimit

137

130

%RPD

2.02

0

**RPDLimit** 

20

0

Qual

%REC

114

99.4

- Е Value above quantitation range
- J
- Analyte detected below quantitation limits
- Page 16 of 32

- P Sample pH Not In Range
- RLReporting Detection Limit
- Sample container temperature is out of limit as specified

## Hall Environmental Analysis Laboratory, Inc.

1.2

WO#: 1901787

06-Feb-19

**Client:** Navajo Refining Company

**Project: RO** Reject

Surr: Tetrachloro-m-xylene

Sample ID MB-42803	SampT	ype: ME	BLK	Tes	tCode: E	PA Method	8082A: PCB's	3		
Client ID: PBW	Batch	n ID: <b>42</b>	803	F	RunNo: <b>5</b>	7368				
Prep Date: 1/24/2019	Analysis D	ate: 1/	30/2019	5	SeqNo: 1	919398	Units: µg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Aroclor 1016	ND	1.0								
Aroclor 1221	ND	1.0								
Aroclor 1232	ND	1.0								
Aroclor 1242	ND	1.0								
Aroclor 1248	ND	1.0								
Aroclor 1254	ND	1.0								
Aroclor 1260	ND	1.0								
Surr: Decachlorobiphenyl	1.9		2.500		76.0	24.8	102			
Surr: Tetrachloro-m-xylene	0.95		2.500		38.0	15.6	106			
Sample ID LCS-42803	SampT	ype: <b>LC</b>	s	Tes	tCode: E	PA Method	8082A: PCB's	5		
Client ID: LCSW	Batch	n ID: <b>42</b>	803	F	RunNo: <b>5</b>	7368				
Prep Date: 1/24/2019	Analysis D	ate: 1/	30/2019	5	SeqNo: 1	919399	Units: µg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Aroclor 1016	2.6	1.0	5.000	0	52.6	25.9	120			
Aroclor 1260	2.7	1.0	5.000	0	54.5	38.4	134			
Surr: Decachlorobiphenyl	1.3		2.500		52.0	24.8	102			

Sample ID LCSD-42803	SampT	ype: <b>LC</b>	SD	Tes	tCode: El	PA Method	8082A: PCB's	6		
Client ID: LCSS02	Batch	ID: <b>42</b>	803	F	RunNo: 5	7368				
Prep Date: 1/24/2019	Analysis D	ate: 1/	30/2019	8	SeqNo: 1	919400	Units: µg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Aroclor 1016	3.9	1.0	5.000	0	77.8	25.9	120	38.5	17.9	R
Aroclor 1260	3.9	1.0	5.000	0	78.1	38.4	134	35.7	16.2	R
Surr: Decachlorobiphenyl	1.8		2.500		74.0	24.8	102	0	0	
Surr: Tetrachloro-m-xylene	1.8		2.500		71.2	15.6	106	0	0	

46.8

15.6

106

2.500

#### Qualifiers:

Value exceeds Maximum Contaminant Level.

Sample Diluted Due to Matrix D

Η Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

Practical Quanitative Limit

% Recovery outside of range due to dilution or matrix

В Analyte detected in the associated Method Blank

E Value above quantitation range

Reporting Detection Limit

J Analyte detected below quantitation limits

P Sample pH Not In Range

RL

Sample container temperature is out of limit as specified

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#### Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID rb2 SampType: MBLK TestCode: EPA Method 8260B: VOLATILES PBW Client ID: Batch ID: **B57169** RunNo: 57169 Prep Date: Analysis Date: 1/22/2019 SeqNo: 1912409 Units: µg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Benzene ND 1.0 ND Toluene 1.0 ND Ethylbenzene 1.0 1,2-Dichloroethane (EDC) ND 1.0 1,2-Dibromoethane (EDB) ND 1.0 Carbon Tetrachloride ND 1.0 Chloroform ND 1.0 ND 1,1-Dichloroethane 1.0 1.1-Dichloroethene ND 1.0 ND Methylene Chloride 3.0 1,1,2,2-Tetrachloroethane ND 2.0 ND Tetrachloroethene (PCE) 1.0 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 Trichloroethene (TCE) ND 1.0 Vinyl chloride ND 1.0 Xylenes, Total ND 1.5 Surr: 1,2-Dichloroethane-d4 10 10.00 103 70 130 Surr: 4-Bromofluorobenzene 10 10.00 101 70 130 Surr: Dibromofluoromethane 10 10.00 101 70 130 Surr: Toluene-d8 11 10.00 108 70 130

Sample ID 100ng lcs2	SampT	ype: <b>LC</b>	s	Tes	tCode: El	PA Method	8260B: VOL	ATILES		
Client ID: LCSW	Batch	n ID: <b>B5</b>	7169	F	RunNo: 5	7169				
Prep Date:	Analysis D	ate: 1/	22/2019	S	SeqNo: 1	912410	Units: µg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Benzene	18	1.0	20.00	0	88.8	70	130			
Toluene	19	1.0	20.00	0	97.2	70	130			
1,1-Dichloroethene	19	1.0	20.00	0	97.4	70	130			
Trichloroethene (TCE)	17	1.0	20.00	0	84.4	70	130			
Surr: 1,2-Dichloroethane-d4	10		10.00		101	70	130			
Surr: 4-Bromofluorobenzene	9.9		10.00		99.3	70	130			
Surr: Dibromofluoromethane	10		10.00		101	70	130			
Surr: Toluene-d8	10		10.00		104	70	130			

#### Qualifiers:

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

in the associated Method Blank

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## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID 1901787-001a ms2 SampType: MS TestCode: EPA Method 8260B: VOLATILES Client ID: R.O. Reject Batch ID: **B57169** RunNo: 57169 Units: µg/L Prep Date: Analysis Date: 1/23/2019 SeqNo: 1912412 Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual 87.9 70 Benzene 18 1.0 20.00 0 130 Toluene 20 1.0 20.00 0 99.3 70 130 95.0 1,1-Dichloroethene 19 1.0 20.00 0 67.6 130 Trichloroethene (TCE) 16 1.0 20.00 0 81.0 70 130 Surr: 1,2-Dichloroethane-d4 11 10.00 108 70 130 Surr: 4-Bromofluorobenzene 10 10.00 99.6 70 130 Surr: Dibromofluoromethane 10 10.00 102 70 130 Surr: Toluene-d8 10 103 70 10.00 130

Sample ID 1901787-001a msd2 SampType: MSD TestCode: EPA Method 8260B: VOLATILES Client ID: R.O. Reject Batch ID: **B57169** RunNo: 57169 Prep Date: Analysis Date: 1/23/2019 SeqNo: 1912413 Units: µg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Benzene 17 1.0 20.00 0 82.6 70 130 6.15 20 Toluene 19 1.0 20.00 0 94.9 70 130 4.58 20 1,1-Dichloroethene 18 20.00 0 92.1 67.6 130 3.08 20 1.0 Trichloroethene (TCE) 16 1.0 20.00 0 79.7 70 130 1.65 20 10.00 105 70 0 Surr: 1,2-Dichloroethane-d4 11 130 Λ Surr: 4-Bromofluorobenzene 10 10.00 101 70 130 0 0 Surr: Dibromofluoromethane 10 102 70 0 0 10.00 130 Surr: Toluene-d8 9.9 10.00 99.2 70 130 0 0

Sample ID 100ng Ics	SampT	ype: LC	s	Tes	tCode: El	PA Method	8260B: VOL	ATILES		
Client ID: LCSW	Batch	n ID: <b>B5</b>	7171	R	RunNo: 5	7171				
Prep Date:	Analysis D	ate: 1/	22/2019	S	SeqNo: 1	912422	Units: µg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Benzene	21	1.0	20.00	0	104	70	130			
Toluene	20	1.0	20.00	0	99.3	70	130			
1,1-Dichloroethene	20	1.0	20.00	0	101	70	130			
Trichloroethene (TCE)	19	1.0	20.00	0	93.1	70	130			
Surr: 1,2-Dichloroethane-d4	10		10.00		105	70	130			
Surr: 4-Bromofluorobenzene	11		10.00		107	70	130			
Surr: Dibromofluoromethane	11		10.00		106	70	130			
Surr: Toluene-d8	10		10.00		103	70	130			

#### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

PQL Practical Quanitative Limit

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

P Sample pH Not In Range

RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

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## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

**Client:** Navajo Refining Company

**Project:** RO Reject

Sample ID rb SampType: MBLK TestCode: EPA Method 8260B: VOLATILES PBW Client ID: Batch ID: **B57171** RunNo: 57171 Prep Date: Analysis Date: 1/22/2019 SeqNo: 1912429 Units: µg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Benzene ND 1.0 ND Toluene 1.0 ND Ethylbenzene 1.0 1,2-Dichloroethane (EDC) ND 1.0 1,2-Dibromoethane (EDB) ND 1.0 Carbon Tetrachloride ND 1.0 Chloroform ND 1.0 1,1-Dichloroethane ND 1.0 1.1-Dichloroethene ND 1.0 ND Methylene Chloride 3.0 1,1,2,2-Tetrachloroethane ND 2.0 Tetrachloroethene (PCE) ND 1.0 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 Trichloroethene (TCE) ND 1.0 Vinyl chloride ND 1.0 Xylenes, Total ND 1.5 Surr: 1,2-Dichloroethane-d4 11 10.00 107 70 130 Surr: 4-Bromofluorobenzene 11 10.00 107 70 130 Surr: Dibromofluoromethane 11 10.00 109 70 130 Surr: Toluene-d8 10 10.00 100 70 130

#### Qualifiers:

- \* Value exceeds Maximum Contaminant Level.
- D Sample Diluted Due to Matrix
- H Holding times for preparation or analysis exceeded
- ND Not Detected at the Reporting Limit
- PQL Practical Quanitative Limit
- S % Recovery outside of range due to dilution or matrix
- B Analyte detected in the associated Method Blank
- E Value above quantitation range

Reporting Detection Limit

- J Analyte detected below quantitation limits
- P Sample pH Not In Range

RL

W Sample container temperature is out of limit as specified

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## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID MB-R57443 SampType: MBLK TestCode: EPA 8270D: Semivolatiles Client ID: PBW Batch ID: **R57443** RunNo: 57443 Prep Date: Analysis Date: 1/25/2019 SeqNo: 1921674 Units: µg/L Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual ND 0.010 Atrazine ND 1,2,4-Trichlorobenzene 0.010 ND 0.010 2,4,6-Trichlorophenol 2,4-Dichlorophenol ND 0.010 2,4-Dimethylphenol ND 0.010 2,4-Dinitrophenol ND 0.010 2,4-Dinitrotoluene ND 0.010 2,6-Dinitrotoluene ND 0.010 2-Chloronaphthalene ND 0.0010 ND 0.010 2-Chlorophenol 2-Nitrophenol ND 0.010 3,3´-Dichlorobenzidine ND 0.010 4,6-Dinitro-2-methylphenol ND 0.010 4-Bromophenyl phenyl ether ND 0.010 4-Chloro-3-methylphenol ND 0.010 4-Chlorophenyl phenyl ether ND 0.010 4-Nitrophenol ND 0.010 Acenaphthene ND 0.0010 Acenaphthylene ND 0.0010 Anthracene ND 0.0010 Benzidine ND 0.010 Benzo(g,h,i)perylene ND 0.0010 Benz(a)anthracene 0.0010 ND Benzo(a)pyrene ND 0.0010 Benzo(b)fluoranthene ND 0.0010 Benzo(k)fluoranthene ND 0.0010 Bis(2-chloroethoxy)methane ND 0.010 Bis(2-chloroethyl)ether ND 0.010 Bis(2-chloroisopropyl)ether ND 0.010 Bis(2-ethylhexyl)phthalate 0.0030 ND 0.0030 Butyl benzyl phthalate ND Chrysene ND 0.0010 Dibenz(a,h)anthracene ND 0.0010 Diethyl phthalate ND 0.0030 ND 0.0030 Dimethyl phthalate Di-n-butyl phthalate ND 0.0030 ND 0.0030 Di-n-octyl phthalate Fluoranthene ND 0.0010 ND 0.0010 Fluorene

#### Qualifiers:

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

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## Hall Environmental Analysis Laboratory, Inc.

SampType: LCS

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID LCS-R57443

Sample ID MB-R57443 SampType: MBLK TestCode: EPA 8270D: Semivolatiles PBW Client ID: Batch ID: **R57443** RunNo: 57443 Analysis Date: 1/25/2019 Prep Date: SeqNo: 1921674 Units: µg/L Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Hexachlorobenzene ND 0.0010 ND Hexachlorobutadiene 0.010 ND Hexachlorocyclopentadiene 0.010 Hexachloroethane ND 0.010 Indeno(1,2,3-cd)pyrene ND 0.0010 0.010 Isophorone ND Naphthalene ND 0.0010 ND 0.010 Nitrobenzene N-Nitrosodimethylamine ND 0.010 N-Nitrosodi-n-propylamine ND 0.010 N-Nitrosodiphenylamine ND 0.010 Pentachlorophenol ND 0.010 0.0010 Phenanthrene ND Phenol ND 0.010 ND 0.0010 Pyrene

Client ID: LCSW	Batch	ID: <b>R57443</b>	F	RunNo: 5	7443				
Prep Date:	Analysis D	ate: 1/25/2019	9	SeqNo: 1	921675	Units: µg/L			
Analyte	Result	PQL SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Atrazine	0.039	0.05000	0	78.2	39	141			
1,2,4-Trichlorobenzene	0.026	0.05000	0	53.0	24	120			
2,4,6-Trichlorophenol	0.037	0.05000	0	73.8	42	120			
2,4-Dichlorophenol	0.029	0.05000	0	58.6	36	120			
2,4-Dimethylphenol	0.029	0.05000	0	58.6	33	120			
2,4-Dinitrophenol	0.043	0.05000	0	85.4	10	120			
2,4-Dinitrotoluene	0.039	0.05000	0	78.6	49	124			
2,6-Dinitrotoluene	0.036	0.05000	0	71.4	46	120			
2-Chloronaphthalene	0.030	0.05000	0	61.0	37	120			
2-Chlorophenol	0.029	0.05000	0	57.2	25	120			
2-Nitrophenol	0.033	0.05000	0	65.6	31	120			
3,3´-Dichlorobenzidine	0.037	0.05000	0	73.8	44	120			
4,6-Dinitro-2-methylphenol	0.042	0.05000	0	84.4	38	138			
4-Bromophenyl phenyl ether	0.037	0.05000	0	74.8	45	120			
4-Chloro-3-methylphenol	0.031	0.05000	0	62.6	40	120			
4-Chlorophenyl phenyl ether	0.034	0.05000	0	68.0	44	120			
4-Nitrophenol	0.015	0.05000	0	29.2	10	120			
Acenaphthene	0.036	0.05000	0	72.4	41	120			

#### Qualifiers:

- \* Value exceeds Maximum Contaminant Level.
- D Sample Diluted Due to Matrix
- H Holding times for preparation or analysis exceeded
- ND Not Detected at the Reporting Limit
- PQL Practical Quanitative Limit
- S % Recovery outside of range due to dilution or matrix
- B Analyte detected in the associated Method Blank

TestCode: EPA 8270D: Semivolatiles

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

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## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

**Client:** Navajo Refining Company

**Project:** RO Reject

Prep Date:   Analysis Date:   1/25/2019   Seq* No:   1921675   Units:   µg/L	Sample ID LCS-R57443	SampT	ype: <b>LC</b>	S	Tes	tCode: El	PA 8270D:	Semivolatiles			
Analyte Result PQL SPK value SPK Ref Val %REC LowLimit High Limit %RPD RPD Limit Qual Acanaphthylene 0.035 0.05000 0 70.4 43 120 120 120 120 120 120 120 120 120 120	Client ID: LCSW	Batch	n ID: R5	57443	R	RunNo: <b>5</b>	7443				
Accenaphthylene         0.035         0.05000         0         70.4         43         120           Anthracene         0.038         0.05000         0         76.8         45         120           Benz/Gla, Diperylene         0.043         0.05000         0         38.6         1         120           Benz/Gla, Diperylene         0.043         0.05000         0         77.0         47         120           Benz/Glayprene         0.040         0.05000         0         77.0         47         120           Benz/Glayprene         0.040         0.05000         0         84.2         46         120           Benz/Cyblitoranthene         0.041         0.05000         0         84.2         46         120           Bis/C-chloroethylpether         0.033         0.05000         0         65.4         23         120           Bis/C-chropethylpether         0.033         0.05000         0         65.4         23         120           Bis/C-chropethylphthalate         0.039         0.05000         0         77.8         43         122           Burly benzyl phthalate         0.039         0.05000         0         77.8         47         120	Prep Date:	Analysis D	ate: 1	/25/2019	S	SeqNo: 1	921675	Units: µg/L			
Anthracene         0.038         0.05000         0         76.8         45         120           Benzidine         0.019         0.05000         0         38.6         1         120           Benzid jahrperine         0.043         0.05000         0         88.8         48         121           Benzid jahracene         0.038         0.05000         0         77.0         47         120           Benzid jahracene         0.040         0.05000         0         79.2         47         120           Benzid jahracene         0.041         0.05000         0         84.2         46         120           Benzid jahracene         0.041         0.05000         0         61.8         33         120           Bis(2-chlorabox)melhane         0.031         0.05000         0         61.8         33         120           Bis(2-chlorabox)melhane         0.032         0.05000         0         61.8         33         120           Bis(2-chlorabox)melhane         0.032         0.05000         0         77.8         43         122           Bis(2-chlorabox)melhane         0.032         0.05000         0         77.8         43         122	Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Benzidine         0.019         0.05000         0         38.6         1         120           Benzidaj hiperylene         0.043         0.05000         0         77.0         47         120           Benzidaj hiperylene         0.040         0.05000         0         77.0         47         120           Benzidyljurarihene         0.042         0.05000         0         84.2         46         120           Bisiz-chloroethorylmehane         0.031         0.05000         0         81.2         46         120           Bisiz-chloroethorylmehane         0.033         0.05000         0         65.4         23         120           Bisiz-chloroethylpeher         0.033         0.05000         0         66.4         23         120           Bisiz-chryphinbalate         0.039         0.05000         0         77.8         43         122           Buly benzyl phihalate         0.039         0.05000         0         77.2         43         121           Chrysene         0.039         0.05000         0         77.8         48         122           Dimetryl phihalate         0.035         0.05000         0         77.6         48         122	Acenaphthylene	0.035		0.05000	0	70.4	43	120			
Benzolgi, h)perylene         0.043         0.05000         0         85.8         48         121           Benzola)hribacene         0.038         0.05000         0         77.0         47         120           Benzola)hyperne         0.040         0.05000         0         79.2         47         120           Benzola)hyperne         0.041         0.05000         0         84.2         46         120           Benzola, hyperhymethane         0.031         0.05000         0         61.8         33         120           Bis(2-chlorostopropylether         0.032         0.05000         0         65.4         23         120           Bis(2-chlorostopropylether         0.032         0.05000         0         64.0         28         120           Bis(2-thlyberdy) phthalate         0.039         0.05000         0         77.2         43         121           Chrysene         0.037         0.05000         0         77.4         48         122           Dimetryla phthalate         0.039         0.05000         0         77.6         48         122           Dimetryla phthalate         0.039         0.05000         0         77.6         48         122	Anthracene	0.038		0.05000	0	76.8	45	120			
Benz(a)anithracene         0.038         0.05000         0         77.0         47         120           Benz(a)pyene         0.040         0.05000         0         79.2         47         120           Benz(a)filturanthene         0.042         0.05000         0         81.2         46         120           Bis(2-chloroethoxylmethane         0.031         0.05000         0         61.8         33         120           Bis(2-chloroethoxylmethane         0.033         0.05000         0         65.4         23         120           Bis(2-chloroethoxylphthalate         0.032         0.05000         0         66.4         28         120           Bis(2-chloroethoxylphthalate         0.039         0.05000         0         77.8         43         122           Bis(2-chlorylphthalate         0.039         0.05000         0         77.8         43         122           Buyl benzyl phthalate         0.039         0.05000         0         77.8         47         120           Diehryl phthalate         0.039         0.05000         0         77.6         48         122           Diendryl phthalate         0.039         0.05000         0         78.6         49	Benzidine	0.019		0.05000	0	38.6	1	120			
Benzo (a) pyrene         0.040         0.05000         0         79.2         47         120           Benzo (b) Huranthene         0.042         0.05000         0         84.2         46         120           Benzo (b) Huranthene         0.041         0.05000         0         81.2         46         120           Bis (2-chlorosbry)methane         0.031         0.05000         0         61.8         33         120           Bis (2-chlorosbryopy)ether         0.032         0.05000         0         64.0         28         120           Bis (2-chlorosbryopy)ether         0.032         0.05000         0         64.0         28         120           Bis (2-chlorosbryopy)ether         0.032         0.05000         0         77.8         43         121           Bis (2-chlorosbryopy)ether         0.032         0.05000         0         77.8         43         122           Bis (2-chlorosbryopy)ether         0.032         0.05000         0         77.4         48         120           Bis (2-chlorosbryopy)ether         0.039         0.05000         0         77.6         48         122           Diber (a) philatale         0.039         0.05000         0         78.6	Benzo(g,h,i)perylene	0.043		0.05000	0	85.8	48	121			
Benzo(h)fluoranthene         0.042         0.05000         0         84.2         46         120           Benzo(h)fluoranthene         0.041         0.05000         0         81.2         46         120           Bis(2-chloroethoy)methane         0.031         0.05000         0         61.8         33         120           Bis(2-chlorostopyplether         0.033         0.05000         0         64.0         28         120           Bis(2-chlorostopyplether         0.032         0.05000         0         64.0         28         120           Bis(2-chlorostopyplether         0.039         0.05000         0         77.2         43         122           Bis(y benzyl phthalate         0.039         0.05000         0         77.8         43         122           Chrysene         0.037         0.05000         0         77.8         47         120           Dibertyl phthalate         0.039         0.05000         0         77.8         48         120           Di-n-byl phthalate         0.035         0.05000         0         78.6         48         120           Di-n-otyl phthalate         0.037         0.05000         0         78.6         49         121<	Benz(a)anthracene	0.038		0.05000	0	77.0	47	120			
Benzo(k)fluoranthene         0.041         0.05000         0         81.2         46         120           Bis(2-chloroethoxy)methane         0.031         0.05000         0         61.8         33         120           Bis(2-chloroethoxy)methane         0.032         0.05000         0         64.0         28         120           Bis(2-chlorostryopy)phther         0.032         0.05000         0         64.0         28         120           Bis(2-chlorostryopy)phthalate         0.039         0.05000         0         77.8         43         122           Buty benzyl phthalate         0.039         0.05000         0         77.4         43         120           Dibenz(a,h)anihracene         0.039         0.05000         0         77.8         47         120           Dibenz(a,h)anihracene         0.039         0.05000         0         77.8         47         120           Dibenz(a,h)anihracene         0.039         0.05000         0         77.6         48         122           Dimethyl phthalate         0.039         0.05000         0         78.6         49         121           Di-butyl phthalate         0.041         0.05000         0         81.8 <th< td=""><td>Benzo(a)pyrene</td><td>0.040</td><td></td><td>0.05000</td><td>0</td><td>79.2</td><td>47</td><td>120</td><td></td><td></td><td></td></th<>	Benzo(a)pyrene	0.040		0.05000	0	79.2	47	120			
Bis(2-chioroethoxy)methane         0.031         0.05000         0         61.8         33         120           Bis(2-chioroethy)lether         0.033         0.05000         0         65.4         23         120           Bis(2-chiorosporpyl)ether         0.032         0.05000         0         64.0         28         120           Bis(2-chiorosporpyl)ether         0.039         0.05000         0         77.2         43         122           Butyl benzyl phthalate         0.039         0.05000         0         77.4         48         120           Dibenz(a,h)anthracene         0.039         0.05000         0         77.8         47         120           Dihendyl phthalate         0.039         0.05000         0         77.6         48         120           Dihendyl phthalate         0.035         0.05000         0         70.4         48         120           Dihendyl phthalate         0.039         0.05000         0         78.6         49         121           Dihendyl phthalate         0.034         0.05000         0         81.8         42         125           Fluoranthene         0.034         0.05000         0         67.6         47	Benzo(b)fluoranthene	0.042		0.05000	0	84.2	46	120			
Bis(2-chloroethy)ether         0.033         0.05000         0         65.4         23         120           Bis(2-chloroisopropylether         0.032         0.05000         0         64.0         28         120           Bis(2-chloroisopropylether         0.039         0.05000         0         77.2         43         121           Chrysene         0.037         0.05000         0         77.4         48         120           Dibenz(a,h)anthracene         0.039         0.05000         0         77.8         47         120           Diethyl phthalate         0.039         0.05000         0         77.6         48         122           Di-n-butyl phthalate         0.035         0.05000         0         77.6         48         122           Di-n-butyl phthalate         0.039         0.05000         0         70.6         48         122           Di-n-butyl phthalate         0.039         0.05000         0         78.6         49         121           Di-n-butyl phthalate         0.031         0.05000         0         73.8         51         120           Fluoranthene         0.034         0.05000         0         73.8         51         120	Benzo(k)fluoranthene	0.041		0.05000	0	81.2	46	120			
Bis(2-chloroisopropylether         0.032         0.05000         0         64.0         28         120           Bis(2-elthylhexyl)phthalate         0.039         0.05000         0         77.8         43         122           Butyl benzyl phthalate         0.039         0.05000         0         77.2         43         121           Chrysene         0.037         0.05000         0         77.4         48         120           Dibenz(a) fighthalate         0.039         0.05000         0         77.6         48         122           Dimethyl phthalate         0.035         0.05000         0         77.6         48         122           Din-butyl phthalate         0.035         0.05000         0         70.4         48         122           Din-butyl phthalate         0.035         0.05000         0         70.4         48         120           Din-butyl phthalate         0.037         0.05000         0         73.8         51         120           Fluoranthene         0.037         0.05000         0         73.8         51         120           Hexachlorobutatiene         0.035         0.05000         0         57.8         19         120 <td>Bis(2-chloroethoxy)methane</td> <td>0.031</td> <td></td> <td>0.05000</td> <td>0</td> <td>61.8</td> <td>33</td> <td>120</td> <td></td> <td></td> <td></td>	Bis(2-chloroethoxy)methane	0.031		0.05000	0	61.8	33	120			
Bis(2-ethylhexyl)phthalate   0.039   0.05000   0   77.8   43   122	Bis(2-chloroethyl)ether	0.033		0.05000	0	65.4	23	120			
Buyl benzyl phthalate         0.039         0.05000         0         77.2         43         121           Chrysene         0.037         0.05000         0         74.4         48         120           Dibenz(a) h)anthracene         0.039         0.05000         0         77.8         47         120           Dienbyl phthalate         0.039         0.05000         0         77.6         48         122           Dimethyl phthalate         0.035         0.05000         0         78.6         49         121           Di-n-butyl phthalate         0.039         0.05000         0         78.6         49         121           Di-n-butyl phthalate         0.039         0.05000         0         78.6         49         121           Di-n-butyl phthalate         0.034         0.05000         0         73.8         51         120           Fluorand         0.037         0.05000         0         67.6         47         120           Hexachlorobutadiene         0.035         0.05000         0         57.8         19         120           Hexachlorocyclopentadiene         0.022         0.05000         0         57.0         15         120	Bis(2-chloroisopropyl)ether	0.032		0.05000	0	64.0	28	120			
Chrysene         0.037         0.05000         0         74.4         48         120           Dibenz(a,h)anthracene         0.039         0.05000         0         77.8         47         120           Diethyl phthalate         0.039         0.05000         0         77.6         48         122           Dir-bulyl phthalate         0.035         0.05000         0         70.4         48         120           Dir-bulyl phthalate         0.039         0.05000         0         78.6         49         121           Dir-bulyl phthalate         0.041         0.05000         0         73.8         51         120           Fluoranthene         0.037         0.05000         0         73.8         51         120           Fluorene         0.034         0.05000         0         67.6         47         120           Hexachlorobutadiene         0.029         0.05000         0         57.8         19         120           Hexachlorocyclopentadiene         0.022         0.05000         0         57.8         19         122           Indean(1,2,3-d)pyrene         0.038         0.05000         0         57.0         49         122 <t< td=""><td>Bis(2-ethylhexyl)phthalate</td><td>0.039</td><td></td><td>0.05000</td><td>0</td><td>77.8</td><td>43</td><td>122</td><td></td><td></td><td></td></t<>	Bis(2-ethylhexyl)phthalate	0.039		0.05000	0	77.8	43	122			
Dibenz(a,h)anthracene   0.039   0.05000   0 77.8   47   120     Diethyl phthalate   0.039   0.05000   0 77.6   48   122     Dimethyl phthalate   0.035   0.05000   0 70.4   48   120     Di-n-bulyl phthalate   0.039   0.05000   0 78.6   49   121     Di-n-bulyl phthalate   0.039   0.05000   0 81.8   42   125     Fluoranthene   0.037   0.05000   0 73.8   51   120     Fluoranthene   0.034   0.05000   0 73.8   51   120     Fluorene   0.034   0.05000   0 67.6   47   120     Hexachlorobenzene   0.035   0.05000   0 70.0   44   120     Hexachlorocyclopentadiene   0.029   0.05000   0 57.8   19   120     Hexachlorocyclopentadiene   0.022   0.05000   0 44.0   15   120     Indeno(1,2,3-cd)pyrene   0.038   0.05000   0 75.0   15   120     Indeno(1,2,3-cd)pyrene   0.038   0.05000   0 75.0   49   122     Isophorone   0.031   0.05000   0 61.2   36   120     Naphthalene   0.027   0.05000   0 61.2   36   120     Naphthalene   0.027   0.05000   0 53.8   27   120     Nitrobenzene   0.030   0.05000   0 59.2   27   120     N-Nitrosodimethylamine   0.019   0.05000   0 64.0   31   120     N-Nitrosodimethylamine   0.032   0.05000   0 64.0   31   120     Pentachlorophenol   0.038   0.05000   0 76.0   23   120     Pentachlorophenol   0.038   0.05000   0 76.0   23   120     Pentanthrene   0.034   0.05000   0 76.0   23   120     Pentanthrene   0.034   0.05000   0 76.0   23   120     Phenanthrene   0.034   0.05000   0 76.0   26.2   10     Phenanthr	Butyl benzyl phthalate	0.039		0.05000	0	77.2	43	121			
Diethyl phthalate         0.039         0.05000         0         77.6         48         122           Dimethyl phthalate         0.035         0.05000         0         70.4         48         120           Di-n-bulyl phthalate         0.039         0.05000         0         78.6         49         121           Di-n-octyl phthalate         0.041         0.05000         0         81.8         42         125           Fluoranthene         0.037         0.05000         0         73.8         51         120           Hexachlorobenzene         0.034         0.05000         0         67.6         47         120           Hexachlorobutadiene         0.035         0.05000         0         57.8         19         120           Hexachlorocyclopentadiene         0.022         0.05000         0         57.8         19         120           Hexachlorocyclopentadiene         0.022         0.05000         0         57.0         15         120           Hexachlorocyclopentadiene         0.028         0.05000         0         57.0         15         120           Hexachlorophoroe         0.031         0.05000         0         61.2         36         120	Chrysene	0.037		0.05000	0	74.4	48	120			
Dimethyl phthalate         0.035         0.05000         0         70.4         48         120           Di-n-butyl phthalate         0.039         0.05000         0         78.6         49         121           Di-n-octyl phthalate         0.041         0.05000         0         81.8         42         125           Fluoranthene         0.037         0.05000         0         67.6         47         120           Hexachlorobenzene         0.035         0.05000         0         67.6         47         120           Hexachlorobutadiene         0.029         0.05000         0         57.8         19         120           Hexachlorocyclopentaldiene         0.022         0.05000         0         57.8         19         120           Hexachlorocyclopentaldiene         0.022         0.05000         0         57.0         15         120           Hexachlorocyclopentaldiene         0.028         0.05000         0         57.0         15         120           Hexachlorocyclopentaldiene         0.028         0.05000         0         57.0         49         122           Isophorone         0.031         0.05000         0         53.8         27         120 <td>Dibenz(a,h)anthracene</td> <td>0.039</td> <td></td> <td>0.05000</td> <td>0</td> <td>77.8</td> <td>47</td> <td>120</td> <td></td> <td></td> <td></td>	Dibenz(a,h)anthracene	0.039		0.05000	0	77.8	47	120			
Di-n-buly phthalate         0.039         0.05000         0         78.6         49         121           Di-n-octyl phthalate         0.041         0.05000         0         81.8         42         125           Fluoranthene         0.037         0.05000         0         73.8         51         120           Fluorene         0.034         0.05000         0         67.6         47         120           Hexachlorobenzene         0.035         0.05000         0         70.0         44         120           Hexachlorocyclopentadiene         0.029         0.05000         0         57.8         19         120           Hexachlorocyclopentadiene         0.022         0.05000         0         44.0         15         120           Hexachlorocyclopentadiene         0.028         0.05000         0         57.0         15         120           Hexachlorocyclopentadiene         0.028         0.05000         0         75.0         49         122           Hexachlorocyclopentadiene         0.038         0.05000         0         61.2         36         120           Hexachlorocyclopentore         0.031         0.05000         0         53.8         27         120	Diethyl phthalate	0.039		0.05000	0	77.6	48	122			
Di-n-otly phthalate         0.041         0.05000         0         81.8         42         125           Fluoranthene         0.037         0.05000         0         73.8         51         120           Fluorene         0.034         0.05000         0         67.6         47         120           Hexachlorobenzene         0.035         0.05000         0         70.0         44         120           Hexachlorobutadiene         0.029         0.05000         0         57.8         19         120           Hexachlorocyclopentadiene         0.022         0.05000         0         57.8         19         120           Hexachlorocethane         0.022         0.05000         0         57.0         15         120           Indeno(1,2,3-cd)pyrene         0.038         0.05000         0         75.0         49         122           Isophorone         0.031         0.05000         0         61.2         36         120           Naphthalene         0.027         0.05000         0         53.8         27         120           N-Nitrosodimethylamine         0.019         0.05000         0         64.0         31         120           N-Ni	Dimethyl phthalate	0.035		0.05000	0	70.4	48	120			
Fluoranthene         0.037         0.05000         0         73.8         51         120           Fluorene         0.034         0.05000         0         67.6         47         120           Hexachlorobenzene         0.035         0.05000         0         70.0         44         120           Hexachlorobutadiene         0.029         0.05000         0         57.8         19         120           Hexachlorocyclopentadiene         0.022         0.05000         0         44.0         15         120           Hexachloroethane         0.028         0.05000         0         57.0         15         120           Indeno(1,2,3-cd)pyrene         0.038         0.05000         0         75.0         49         122           Isophorone         0.031         0.05000         0         61.2         36         120           Naphthalene         0.027         0.05000         0         53.8         27         120           N-Nitrosodimethylamine         0.019         0.05000         0         38.6         10         120           N-Nitrosodiphenylamine         0.032         0.05000         0         65.0         47         120           Pe	Di-n-butyl phthalate	0.039		0.05000	0	78.6	49	121			
Fluorene         0.034         0.05000         0         67.6         47         120           Hexachlorobenzene         0.035         0.05000         0         70.0         44         120           Hexachlorobutadiene         0.029         0.05000         0         57.8         19         120           Hexachlorocyclopentadiene         0.022         0.05000         0         44.0         15         120           Hexachloroethane         0.028         0.05000         0         57.0         15         120           Indeno(1,2,3-cd)pyrene         0.038         0.05000         0         75.0         49         122           Isophorone         0.031         0.05000         0         61.2         36         120           Naphthalene         0.027         0.05000         0         53.8         27         120           Nitrobenzene         0.030         0.05000         0         38.6         10         120           N-Nitrosodir-propylamine         0.032         0.05000         0         64.0         31         120           N-Nitrosodiphenylamine         0.038         0.05000         0         65.0         47         120	Di-n-octyl phthalate	0.041		0.05000	0	81.8	42	125			
Hexachlorobenzene         0.035         0.05000         0         70.0         44         120           Hexachlorobutadiene         0.029         0.05000         0         57.8         19         120           Hexachlorocyclopentadiene         0.022         0.05000         0         44.0         15         120           Hexachloroethane         0.028         0.05000         0         57.0         15         120           Indeno(1,2,3-cd)pyrene         0.038         0.05000         0         75.0         49         122           Isophorone         0.031         0.05000         0         61.2         36         120           Naphthalene         0.027         0.05000         0         53.8         27         120           Nitrobenzene         0.030         0.05000         0         59.2         27         120           N-Nitrosodimethylamine         0.019         0.05000         0         64.0         31         120           N-Nitrosodiphenylamine         0.032         0.05000         0         65.0         47         120           Pentachlorophenol         0.038         0.05000         0         67.0         46         120	Fluoranthene	0.037		0.05000	0	73.8	51	120			
Hexachlorobutadiene       0.029       0.05000       0       57.8       19       120         Hexachlorocyclopentadiene       0.022       0.05000       0       44.0       15       120         Hexachloroethane       0.028       0.05000       0       57.0       15       120         Indeno(1,2,3-cd)pyrene       0.038       0.05000       0       75.0       49       122         Isophorone       0.031       0.05000       0       61.2       36       120         Naphthalene       0.027       0.05000       0       53.8       27       120         Nitrobenzene       0.030       0.05000       0       59.2       27       120         N-Nitrosodimethylamine       0.019       0.05000       0       38.6       10       120         N-Nitrosodiphenylamine       0.032       0.05000       0       64.0       31       120         Pentachlorophenol       0.038       0.05000       0       76.0       23       120         Phenanthrene       0.034       0.05000       0       67.0       46       120         Phenol       0.013       0.05000       0       26.2       10       120 </td <td>Fluorene</td> <td>0.034</td> <td></td> <td>0.05000</td> <td>0</td> <td>67.6</td> <td>47</td> <td>120</td> <td></td> <td></td> <td></td>	Fluorene	0.034		0.05000	0	67.6	47	120			
Hexachlorocyclopentadiene       0.022       0.05000       0       44.0       15       120         Hexachloroethane       0.028       0.05000       0       57.0       15       120         Indeno(1,2,3-cd)pyrene       0.038       0.05000       0       75.0       49       122         Isophorone       0.031       0.05000       0       61.2       36       120         Naphthalene       0.027       0.05000       0       53.8       27       120         Nitrobenzene       0.030       0.05000       0       59.2       27       120         N-Nitrosodimethylamine       0.019       0.05000       0       38.6       10       120         N-Nitrosodiphenylamine       0.032       0.05000       0       64.0       31       120         N-Nitrosodiphenylamine       0.032       0.05000       0       65.0       47       120         Pentachlorophenol       0.038       0.05000       0       76.0       23       120         Phenanthrene       0.034       0.05000       0       67.0       46       120         Phenol       0.013       0.05000       0       26.2       10       120	Hexachlorobenzene	0.035		0.05000	0	70.0	44	120			
Hexachloroethane         0.028         0.05000         0         57.0         15         120           Indeno(1,2,3-cd)pyrene         0.038         0.05000         0         75.0         49         122           Isophorone         0.031         0.05000         0         61.2         36         120           Naphthalene         0.027         0.05000         0         53.8         27         120           Nitrobenzene         0.030         0.05000         0         59.2         27         120           N-Nitrosodimethylamine         0.019         0.05000         0         38.6         10         120           N-Nitrosodiphenylamine         0.032         0.05000         0         64.0         31         120           Pentachlorophenol         0.038         0.05000         0         65.0         47         120           Phenanthrene         0.034         0.05000         0         67.0         23         120           Phenol         0.013         0.05000         0         67.0         46         120	Hexachlorobutadiene	0.029		0.05000	0	57.8	19	120			
Indeno(1,2,3-cd)pyrene         0.038         0.05000         0         75.0         49         122           Isophorone         0.031         0.05000         0         61.2         36         120           Naphthalene         0.027         0.05000         0         53.8         27         120           Nitrobenzene         0.030         0.05000         0         59.2         27         120           N-Nitrosodimethylamine         0.019         0.05000         0         38.6         10         120           N-Nitrosodiphenylamine         0.032         0.05000         0         64.0         31         120           Pentachlorophenol         0.032         0.05000         0         65.0         47         120           Phenanthrene         0.034         0.05000         0         76.0         23         120           Phenol         0.013         0.05000         0         67.0         46         120	Hexachlorocyclopentadiene	0.022		0.05000	0	44.0	15	120			
Isophorone         0.031         0.05000         0         61.2         36         120           Naphthalene         0.027         0.05000         0         53.8         27         120           Nitrobenzene         0.030         0.05000         0         59.2         27         120           N-Nitrosodimethylamine         0.019         0.05000         0         38.6         10         120           N-Nitrosodiphenylamine         0.032         0.05000         0         64.0         31         120           N-Nitrosodiphenylamine         0.032         0.05000         0         65.0         47         120           Pentachlorophenol         0.038         0.05000         0         76.0         23         120           Phenanthrene         0.034         0.05000         0         67.0         46         120           Phenol         0.013         0.05000         0         26.2         10         120	Hexachloroethane	0.028		0.05000	0	57.0	15	120			
Naphthalene       0.027       0.05000       0       53.8       27       120         Nitrobenzene       0.030       0.05000       0       59.2       27       120         N-Nitrosodimethylamine       0.019       0.05000       0       38.6       10       120         N-Nitrosodiphenylamine       0.032       0.05000       0       64.0       31       120         N-Nitrosodiphenylamine       0.032       0.05000       0       65.0       47       120         Pentachlorophenol       0.038       0.05000       0       76.0       23       120         Phenanthrene       0.034       0.05000       0       67.0       46       120         Phenol       0.013       0.05000       0       26.2       10       120	Indeno(1,2,3-cd)pyrene	0.038		0.05000	0	75.0	49	122			
Nitrobenzene         0.030         0.05000         0         59.2         27         120           N-Nitrosodimethylamine         0.019         0.05000         0         38.6         10         120           N-Nitrosodiphenylamine         0.032         0.05000         0         64.0         31         120           N-Nitrosodiphenylamine         0.032         0.05000         0         65.0         47         120           Pentachlorophenol         0.038         0.05000         0         76.0         23         120           Phenanthrene         0.034         0.05000         0         67.0         46         120           Phenol         0.013         0.05000         0         26.2         10         120	Isophorone	0.031		0.05000	0	61.2	36	120			
N-Nitrosodimethylamine       0.019       0.05000       0       38.6       10       120         N-Nitrosodi-n-propylamine       0.032       0.05000       0       64.0       31       120         N-Nitrosodiphenylamine       0.032       0.05000       0       65.0       47       120         Pentachlorophenol       0.038       0.05000       0       76.0       23       120         Phenanthrene       0.034       0.05000       0       67.0       46       120         Phenol       0.013       0.05000       0       26.2       10       120	Naphthalene	0.027		0.05000	0	53.8	27	120			
N-Nitrosodi-n-propylamine       0.032       0.05000       0       64.0       31       120         N-Nitrosodiphenylamine       0.032       0.05000       0       65.0       47       120         Pentachlorophenol       0.038       0.05000       0       76.0       23       120         Phenanthrene       0.034       0.05000       0       67.0       46       120         Phenol       0.013       0.05000       0       26.2       10       120	Nitrobenzene	0.030		0.05000	0	59.2	27	120			
N-Nitrosodiphenylamine       0.032       0.05000       0       65.0       47       120         Pentachlorophenol       0.038       0.05000       0       76.0       23       120         Phenanthrene       0.034       0.05000       0       67.0       46       120         Phenol       0.013       0.05000       0       26.2       10       120	N-Nitrosodimethylamine	0.019		0.05000	0	38.6	10	120			
Pentachlorophenol         0.038         0.05000         0         76.0         23         120           Phenanthrene         0.034         0.05000         0         67.0         46         120           Phenol         0.013         0.05000         0         26.2         10         120	N-Nitrosodi-n-propylamine	0.032		0.05000	0	64.0	31	120			
Phenanthrene         0.034         0.05000         0         67.0         46         120           Phenol         0.013         0.05000         0         26.2         10         120	N-Nitrosodiphenylamine	0.032		0.05000	0	65.0	47	120			
Phenol 0.013 0.05000 0 26.2 10 120	Pentachlorophenol	0.038		0.05000	0	76.0	23	120			
	Phenanthrene	0.034		0.05000	0	67.0	46	120			
Pyrene 0.037 0.05000 0 73.2 47 120	Phenol	0.013		0.05000	0	26.2	10	120			
	Pyrene	0.037		0.05000	0	73.2	47	120			

#### Qualifiers:

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

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## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID MB-42802 SampType: MBLK TestCode: EPA Method 8310: PAHs Client ID: PBW Batch ID: 42802 RunNo: 57348 Prep Date: 1/24/2019 Analysis Date: 1/30/2019 SeqNo: 1919614 Units: µg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Naphthalene ND 3.0 1-Methylnaphthalene ND 3.0 2-Methylnaphthalene ND 3.0 Acenaphthylene ND 3.0 Acenaphthene ND 3.0 Fluorene ND 0.80 Phenanthrene ND 0.60 ND 0.60 Anthracene Fluoranthene ND 0.30 ND Pyrene 0.40 Benz(a)anthracene ND 0.070 ND 0.20 Chrysene Benzo(b)fluoranthene ND 0.10 Benzo(k)fluoranthene ND 0.070 Benzo(a)pyrene ND 0.070 Dibenz(a,h)anthracene ND 0.12 0.12 Benzo(g,h,i)perylene ND Indeno(1,2,3-cd)pyrene ND 0.25 67.1 Surr: Benzo(e)pyrene 13 20.00 48.8 93.3

Sample ID LCS-42802	Samp1	ype: <b>LC</b>	S	Tes	tCode: El	PA Method	8310: PAHs			
Client ID: LCSW	Batcl	n ID: 42	802	R	RunNo: 5	7348				
Prep Date: 1/24/2019	Analysis D	oate: 1/	30/2019	S	SeqNo: 1	919615	Units: µg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Naphthalene	45	3.0	80.00	0	56.7	23.8	80.3			
1-Methylnaphthalene	45	3.0	80.20	0	55.7	23.4	81.9			
2-Methylnaphthalene	45	3.0	80.00	0	56.0	22.9	81.4			
Acenaphthylene	51	3.0	80.20	0	64.0	42.6	86.6			
Acenaphthene	47	3.0	80.00	0	58.6	40.2	83.4			
Fluorene	4.9	0.80	8.020	0	61.6	44.3	85			
Phenanthrene	2.5	0.60	4.020	0	61.4	42	95.2			
Anthracene	2.6	0.60	4.020	0	65.7	57	87.4			
Fluoranthene	5.3	0.30	8.020	0	66.2	55.7	88.9			
Pyrene	4.6	0.40	8.020	0	58.0	49.5	95			
Benz(a)anthracene	0.54	0.070	0.8020	0	67.3	51.9	98.9			
Chrysene	2.6	0.20	4.020	0	63.4	51	95.6			
Benzo(b)fluoranthene	0.64	0.10	1.002	0	63.9	50	95.2			
Benzo(k)fluoranthene	0.33	0.070	0.5000	0	66.0	55.7	91.5			

#### Qualifiers:

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

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## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID LCS-42802 SampType: LCS TestCode: EPA Method 8310: PAHs Client ID: LCSW Batch ID: 42802 RunNo: 57348 Prep Date: 1/24/2019 Analysis Date: 1/30/2019 SeqNo: 1919615 Units: µg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Benzo(a)pyrene 0.32 0.070 0.5020 0 63.7 47.3 98.2 0.68 0 67.9 51.8 Dibenz(a,h)anthracene 0.12 1.002 99.1 0.65 0.12 0 65.0 Benzo(g,h,i)perylene 1.000 51 99.3 Indeno(1,2,3-cd)pyrene 1.3 0.25 2.004 0 66.4 51.5 96.4 Surr: Benzo(e)pyrene 14 20.00 71.1 48.8 93.3

Sample ID LCSD-42802	Samp	Гуре: <b>LC</b>	SD	Tes	tCode: El	PA Method	8310: PAHs			
Client ID: LCSS02	Batc	h ID: <b>42</b> 8	802	F	RunNo: 5	7348				
Prep Date: 1/24/2019	Analysis [	Date: 1/	30/2019	S	SeqNo: 1	919616	Units: µg/L			
Analyte	Result	PQL	SPK value	SPK Ref Val	%REC	LowLimit	HighLimit	%RPD	RPDLimit	Qual
Naphthalene	37	3.0	80.00	0	46.4	23.8	80.3	20.1	34.8	
1-Methylnaphthalene	38	3.0	80.20	0	47.7	23.4	81.9	15.6	33	
2-Methylnaphthalene	38	3.0	80.00	0	47.5	22.9	81.4	16.4	33.3	
Acenaphthylene	45	3.0	80.20	0	56.4	42.6	86.6	12.7	30	
Acenaphthene	42	3.0	80.00	0	52.4	40.2	83.4	11.3	30	
Fluorene	4.5	0.80	8.020	0	56.0	44.3	85	9.54	24.8	
Phenanthrene	2.3	0.60	4.020	0	56.2	42	95.2	8.88	30.2	
Anthracene	2.4	0.60	4.020	0	59.7	57	87.4	9.52	22.3	
Fluoranthene	4.9	0.30	8.020	0	61.3	55.7	88.9	7.62	24.2	
Pyrene	4.3	0.40	8.020	0	53.9	49.5	95	7.36	24.4	
Benz(a)anthracene	0.49	0.070	0.8020	0	61.1	51.9	98.9	9.71	31.3	
Chrysene	2.4	0.20	4.020	0	59.2	51	95.6	6.90	25.5	
Benzo(b)fluoranthene	0.60	0.10	1.002	0	59.9	50	95.2	6.45	25	
Benzo(k)fluoranthene	0.30	0.070	0.5000	0	60.0	55.7	91.5	9.52	32.7	
Benzo(a)pyrene	0.29	0.070	0.5020	0	57.8	47.3	98.2	9.84	33.2	
Dibenz(a,h)anthracene	0.65	0.12	1.002	0	64.9	51.8	99.1	4.51	25.1	
Benzo(g,h,i)perylene	0.61	0.12	1.000	0	61.0	51	99.3	6.35	31.8	
Indeno(1,2,3-cd)pyrene	1.2	0.25	2.004	0	62.4	51.5	96.4	6.20	26.8	
Surr: Benzo(e)pyrene	13		20.00		63.8	48.8	93.3	0		

#### Qualifiers:

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

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#### Hall Environmental Analysis Laboratory, Inc.

WO#: 1901787

06-Feb-19

**Client:** Navajo Refining Company

**Project: RO** Reject

Sample ID MB-42880 SampType: MBLK TestCode: Total Phenolics by SW-846 9067

Client ID: PBW Batch ID: 42880 RunNo: 57339

Prep Date: 1/30/2019 Analysis Date: 1/30/2019 SeqNo: 1918540 Units: µg/L

Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual

Phenolics ND 2.5

Sample ID LCS-42880 SampType: LCS TestCode: Total Phenolics by SW-846 9067

Client ID: LCSW Batch ID: 42880 RunNo: 57339

Prep Date: 1/30/2019 Analysis Date: 1/30/2019 SeqNo: 1918541 Units: µg/L

SPK value SPK Ref Val %REC %RPD **RPDLimit** Analyte Result PQL LowLimit HighLimit Qual

Phenolics 17 2.5 20.00 0 85.6 57.7 149

Sample ID 1901787-001CMS SampType: MS TestCode: Total Phenolics by SW-846 9067

Client ID: R.O. Reject Batch ID: 42880 RunNo: 57339

Prep Date: 1/30/2019 Analysis Date: 1/30/2019 SeqNo: 1918543 Units: µg/L

SPK value SPK Ref Val %REC Result **PQL** LowLimit HighLimit %RPD **RPDLimit** Analyte Qual

Phenolics 20.00 S

Sample ID 1901787-001CMSD SampType: MSD TestCode: Total Phenolics by SW-846 9067

Client ID: Batch ID: 42880 RunNo: 57339 R.O. Reject

Analysis Date: 1/30/2019 Prep Date: 1/30/2019 SeqNo: 1918544 Units: µg/L

Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual

23.8 Phenolics 24 2.5 20.00 0 122 70.1 127 14.3

#### Qualifiers:

Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

Η Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

POL Practical Quanitative Limit

% Recovery outside of range due to dilution or matrix

В Analyte detected in the associated Method Blank

Е Value above quantitation range

J Analyte detected below quantitation limits

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P Sample pH Not In Range

RLReporting Detection Limit

Sample container temperature is out of limit as specified

## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

**Client:** Navajo Refining Company

**Project:** RO Reject

Sample ID MB-R57443 SampType: MBLK TestCode: EPA 335.4: Total Cyanide Subbed

Client ID: PBW Batch ID: R57443 RunNo: 57443

Prep Date: Analysis Date: 1/29/2019 SeqNo: 1921729 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Cyanide ND 0.00500

Sample ID LCS-R57443 SampType: LCS TestCode: EPA 335.4: Total Cyanide Subbed

Client ID: LCSW Batch ID: R57443 RunNo: 57443

Prep Date: Analysis Date: 1/29/2019 SeqNo: 1921730 Units: mg/L

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Cyanide 0.0973 0.1000 0 97.3 85 115

#### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

PQL Practical Quanitative Limit

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

G 1 HN / I D

P Sample pH Not In Range RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

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## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

**Client:** Navajo Refining Company

**Project:** RO Reject

Sample ID Ics-1 99.0uS eC SampType: Ics TestCode: SM2510B: Specific Conductance

Client ID: LCSW Batch ID: R57160 RunNo: 57160

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1911988 Units: µmhos/cm

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Conductivity 98 5.0 99.00 0 98.9 80 120

Sample ID 1901628-002c dup SampType: dup TestCode: SM2510B: Specific Conductance

Client ID: BatchQC Batch ID: R57160 RunNo: 57160

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1911991 Units: µmhos/cm

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Conductivity 6100 5.0 0.394 20

Sample ID 1901748-002c dup SampType: dup TestCode: SM2510B: Specific Conductance

Client ID: BatchQC Batch ID: R57160 RunNo: 57160

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1912002 Units: µmhos/cm

Analyte Result PQL SPK value SPK Ref Val %REC LowLimit HighLimit %RPD RPDLimit Qual

Conductivity 6000 5.0 0.299 20

#### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

PQL Practical Quanitative Limit

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

Reporting Detection Limit

J Analyte detected below quantitation limits

D C 1 HN / I D

P Sample pH Not In Range

RL

W Sample container temperature is out of limit as specified

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## Hall Environmental Analysis Laboratory, Inc.

WO#: **1901787** 

06-Feb-19

Client: Navajo Refining Company

**Project:** RO Reject

Sample ID 1901789-002ams SampType: MS TestCode: EPA Method 8015D: Gasoline Range Client ID: **BatchQC** Batch ID: **R57171** RunNo: 57171 Prep Date: Analysis Date: 1/22/2019 SeqNo: 1912400 Units: mg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual 0.050 O 63.4 0.52 0.5000 104 130

 Gasoline Range Organics (GRO)
 0.52
 0.050
 0.5000
 0
 104
 63.4
 130

 Surr: BFB
 9.8
 10.00
 97.7
 70
 130

Sample ID 1901789-002amsd SampType: MSD TestCode: EPA Method 8015D: Gasoline Range Client ID: **BatchQC** Batch ID: **R57171** RunNo: 57171 Prep Date: Analysis Date: 1/22/2019 SeqNo: 1912401 Units: mg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Gasoline Range Organics (GRO) 0.49 0.050 0.5000 98.2 63.4 130 5.62 20 Surr: BFB 9.7 10.00 96.8 70 130 0

Sample ID 2.5ug gro Ics SampType: LCS TestCode: EPA Method 8015D: Gasoline Range Client ID: LCSW Batch ID: **R57171** RunNo: 57171 Prep Date: Analysis Date: 1/22/2019 SeqNo: 1912406 Units: mg/L SPK value SPK Ref Val %REC %RPD **RPDLimit** Analyte Result **PQL** LowLimit HighLimit Qual Gasoline Range Organics (GRO) 0.53 0.050 0.5000 0 106 70 130 Surr: BFB 70 9.8 10.00 98.0 130

Sample ID rb SampType: MBLK TestCode: EPA Method 8015D: Gasoline Range Client ID: PBW Batch ID: **R57171** RunNo: 57171 Prep Date: Analysis Date: 1/22/2019 SeqNo: 1912407 Units: mg/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Gasoline Range Organics (GRO) ND 0.050 Surr: BFB 9.7 10.00 96.6 70 130

#### Qualifiers:

\* Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

H Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

PQL Practical Quanitative Limit

S % Recovery outside of range due to dilution or matrix

B Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

P Sample pH Not In Range

RL Reporting Detection Limit

W Sample container temperature is out of limit as specified

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## Hall Environmental Analysis Laboratory, Inc.

WO#: 1901787

06-Feb-19

**Client:** Navajo Refining Company

**Project: RO** Reject

Sample ID 1901628-002c dup SampType: dup TestCode: SM4500-H+B / 9040C: pH

Client ID: **BatchQC** Batch ID: **R57160** RunNo: 57160

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1912023 Units: pH units

Analyte Result SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual

7.85 Н рΗ

Sample ID 1901748-002c dup SampType: dup TestCode: SM4500-H+B / 9040C: pH

Client ID: BatchQC Batch ID: **R57160** RunNo: 57160

Prep Date: Analysis Date: 1/21/2019 SeqNo: 1912043 Units: pH units

Analyte Result SPK value SPK Ref Val %REC LowLimit %RPD **RPDLimit** PQL HighLimit Qual

\*H рΗ 8.53

#### Qualifiers:

Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

Η Holding times for preparation or analysis exceeded

ND Not Detected at the Reporting Limit

POL Practical Quanitative Limit

% Recovery outside of range due to dilution or matrix

В Analyte detected in the associated Method Blank

E Value above quantitation range

J Analyte detected below quantitation limits

Page 30 of 32

P Sample pH Not In Range

RL Reporting Detection Limit

Sample container temperature is out of limit as specified

## Hall Environmental Analysis Laboratory, Inc.

0.324

0.555

WO#: 1901787

06-Feb-19

**Client:** Navajo Refining Company

**Project: RO** Reject

Radium-228 ±

Sample ID MB-R57324 SampType: MBLK TestCode: EPA 903.1: Ra 226 and EPA 904.0: Ra 228-Subbed Client ID: PBW Batch ID: **R57324** RunNo: 57324 Prep Date: Analysis Date: 1/29/2019 SeqNo: 1917743 Units: pCi/L Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual Radium-226 0.654 0.754 Radium-226 ± 0.561 0.754 Radium-228 0.636 0.555

#### Qualifiers:

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

Page 31 of 32

## Hall Environmental Analysis Laboratory, Inc.

WO#: 1901787

Qual

06-Feb-19

**Client:** Navajo Refining Company

**Project: RO** Reject

Sample ID MB-42739 SampType: MBLK TestCode: SM2540C MOD: Total Dissolved Solids

Client ID: PBW Batch ID: 42739 RunNo: 57198

Prep Date: Analysis Date: 1/23/2019 SeqNo: 1913205 1/22/2019 Units: mg/L

Analyte Result **PQL** SPK value SPK Ref Val %REC LowLimit HighLimit %RPD **RPDLimit** Qual

Total Dissolved Solids ND 20.0

TestCode: SM2540C MOD: Total Dissolved Solids Sample ID LCS-42739 SampType: LCS

Client ID: LCSW Batch ID: 42739 RunNo: 57198

Units: mg/L Prep Date: 1/22/2019 Analysis Date: 1/23/2019 SeqNo: 1913206

SPK value SPK Ref Val %REC LowLimit %RPD **RPDLimit** Analyte Result **PQL** HighLimit

Total Dissolved Solids 1010 20.0 1000 0 101 120

Sample ID 1901741-003ADUP SampType: DUP TestCode: SM2540C MOD: Total Dissolved Solids

Client ID: **BatchQC** Batch ID: 42739 RunNo: 57198

Prep Date: Analysis Date: 1/23/2019 SeqNo: 1913209 Units: mg/L 1/22/2019

SPK value SPK Ref Val %REC LowLimit Result **PQL** HighLimit %RPD **RPDLimit** Qual Analyte

Total Dissolved Solids 0.608

Sample ID 1901760-002ADUP SampType: DUP TestCode: SM2540C MOD: Total Dissolved Solids

Client ID: Batch ID: 42739 RunNo: 57198 **BatchQC** 

Prep Date: Analysis Date: 1/23/2019 Units: mg/L 1/22/2019 SeqNo: 1913213

SPK value SPK Ref Val %REC LowLimit Analyte Result **PQL** HighLimit %RPD **RPDLimit** Qual

Total Dissolved Solids 578 20.0 0.173 5

#### Qualifiers:

Value exceeds Maximum Contaminant Level.

D Sample Diluted Due to Matrix

Holding times for preparation or analysis exceeded Η

ND Not Detected at the Reporting Limit

POL Practical Quanitative Limit

% Recovery outside of range due to dilution or matrix

В Analyte detected in the associated Method Blank

Е Value above quantitation range

J Analyte detected below quantitation limits

Page 32 of 32

P Sample pH Not In Range

RLReporting Detection Limit

Sample container temperature is out of limit as specified



Hall Environmental Analysis Laboratory
4901 Hawkins NE
Albuquerque, NM 87109

TEL: 505-345-3975 FAX: 505-345-4107 Website: www.hallenvironmental.com

# Sample Log-In Check List

s V dex s V	No   No   No   No   No   No   No   No	Not Present   NA   NA   NA	
s V	No	Not Present ☐	
s V	No	Not Present ☐	
s V	No	NA 🗆	
s V	No	NA 🗆	
s V	No	NA 🗆	
s V	No	NA 🗆	
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	No 🗆	NA 🗀	
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			-1
Date :	Signed By		
		No V  No No No No V  No N	No N

S	hain	-of-Cu	Chain-of-Custody Record	Turn-Around Time:	Time:				I		ū	HALL FNYTDONMENTAL	C	Ž	Ī	F	<	
Client:	Navajo	Client: Navajo Refinery		X Standard	☐ Rush				4		7	ANALYSTS LABORATORY	2 3		<u> </u>			
				Project Name					;	, <u>2</u>	llenvi	www.hallenvironmental.com	ntal c	E		) :		
Mailing	Address	s: P.O. Bo	Mailing Address: P.O. Box 159 Artesia,	R.O. Reject			4	4901 Hawkins NE	awkin	s RE	Ā	- Albuquerane, NM 87109	E Z	M 87	109			
NM 882	NM 88211-0159	6		Project #: P.C	P.O. # 231642		•	Tel. 505-345-3975	5-345	-3975	Ŀ	Fax 505-345-4107	5-345	410				
Phone a	Phone #: 575-748-3311	48-3311									¹nal	Analysis Request	dnes	Į.				
email o	r Fax#: ₹	email or Fax#: 575-746-5451	451	Project Manager:	ger:							(0.						
QA/QC	QA/QC Package:			Scott Denton								<del>7</del> 06						
X Standard	dard		☐ Level 4 (Full Validation)	Robert Combs	S							+1.8			•			
□ Other	<u>*</u>			Sampler: Brad	ad, Historik	K			əpi						-	(1		
	EDD (Type)			On Ice: 💃	X Yes	en e					•			•		7 10		
				Sample Temp	Sample Temperature [ . 3 + 1. 등 (C	5(CF) ≥ 2.8 <sup>©</sup>	_							•		٧)		
Date	Time	Matrix	Sample Request ID	Container Type and #	Preservative Type	HEAL NO. 1901787	8250C: WQC	6010B: WQ	335.4: Tota	7470: Mero 8015: GRO	8082: PCB	Radioactivit Cation/Anio	Phenols SV	Hq	804.1:EDB	eəlddu <b>B</b> 1iA		
6.18	1:32	liquid	R.O. Reject	2 - 500ml P	1-unpres 1- H2SO4	100						×		×		<b>&gt;</b>		
1-18-19 1:35	1:32	liquid	R.O. Reject	3-40ml VOA	HCL		×											
1-18-19	1:35	liquid	R.O. Reject	1-500ml P	HNO3	1			Ĥ	×								
1.18.19 1.35	1:35	liquid	R.O. Reject	1-125ml P	HNO3			×							<del>                                     </del>		_	
1.18.19	(:33	liquid	R.O. Reject	1-500ml P	NaOH				×									
1.18.19	1:35	liquid	R.O. Reject	2-1L P	HNO3							×						
61.81.1	1:35	liquid	R.O. Reject	3-40ml VOA	Na2S2O3										×			
1.18.19 1:35	1:35	pinbil	R.O. Reject	2 - 1L Glass	unpres						×							
61.81.1	1:35	liquid	R.O. Reject	1 - 1L Glass	nnpres		<u>×</u>											
1-18-19	1.35	liquid	R.O. Reject	3-40ml VOA	НСІ					×								
61.81.1	1:35	liquid	R.O. Reject	1-250mlGlas unpres	unpres					×					_			
61.31.1	56:1	liquid	Reject	1 - 1L Glass	H2SO4	T							×	_	_	<u> </u>		
6/31-1	ХХ	liquid	liquid Trip Blank	2-40ml VOA	뙤	200	H					Н						П
Date: Time: 1.18.19 4:00		Relinquishe	Historia (	Received by:	Fedex Lang	Date Time 7 1/2/1	Remarks: email to Scott Denton, Robert Combs and Randy Dade <b>PH2 Trop</b> Metals: As, Al, Ba, B, Cd, Cr, Co, Cu, Fe, Pb, Mn, Hg, Mo, Ni, Se, Ag, U, Zn OCS: 1,1,1-Trichloroethane; 1,1,2,2-Tetrachloroethane; 1,1,2-Tetrachloroethane; 1,1,2-Tetrachloroe	s. email s, Al, Ba, 1,1-Trich	to Scot B, Cd, C loroethar	Dento r, Co, C ie, 1,1,2	n, Robe 1, Fe, Pt 2-Tetrac	rt Com , Mn, Hg hloroeth	os and Mo, Ni, ane: 1,1,	Randy Se, Ag 2,2-Tet	r Dade I, U, Zn rrachlor	oethylene	5.7% ne: 1,12	<b>4</b> 0 %
Date:	Time:	Refinquished by	sd by:	Received by:		Date Time	Inchloroethane; 1,1,2-Irichloroethylene; 1,1-Dichloroethane; 1,1-Dichloroethene; 1,2- Dibromoethane; 1,2-Dichloroethane; Benzene; Carbon Tetrachloride; Chloroform; Dichloromethane; Ethylbenzene; Toluene; Tofal Xylenes; Vinyl Chloride SVOCs: benzo(a)pyrene, phenol, 1-methylnaphthalene, 2-methylnaphthalene, naphthalene	thane; 1, thane; 1,; ethane; P penzo(a);	1,2-Trich 2-Dichlor Ethylben: syrene, p	loroethy oethane æne; To henol, 1	ene; 1,1 Benzer luene; T	-Dichlord e; Carbo stal Xyler aphthale	ethane; n Tetrac ies; Vin) ne, 2-me	1,1-Dic shloride /I Chlori sthylnap	hloroet Chloro ide htthaler	hene; 1, oform; ne, napt	,2- nthalene	
	lf necessary	, samples subr	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	ontracted to other ac	credited laboratories. T	his serves as notice of this	possibility	. Any su	b-contra	ted data	will be	learly no	tated or	the an	alytical	report.		



# Attachment A 2019 Annual Groundwater Monitoring Report, February 2020 (Separate Electronic File)