Final Third Quarter 2013 - Quarterly Groundwater Monitoring Report Outside Tunnel Wells

Red Hill Bulk Fuel Storage Facility Joint Base Pearl Harbor-Hickam, Oahu, Hawaii

DOH Facility ID: 9-102271 DOH Release ID: 990051, 010011, and 020028

September 2013

Department of the Navy Naval Facilities Engineering Command, Hawaii 400 Marshall Road JBPHH, HI 96860-3139



Contract Number N62742-12-D-1853, CTO 0002

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Prepared for:



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Prepared under:

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FINAL THIRD QUARTER 2013 - QUARTERLY GROUNDWATER MONITORING REPORT OUTSIDE TUNNEL WELLS RED HILL BULK FUEL STORAGE FACILITY

Long-Term Groundwater and Soil Vapor Monitoring Red Hill Bulk Fuel Storage Facility Joint-Base Pearl Harbor-Hickam, Oahu, Hawaii

Prepared for:

Department of the Navy Commanding Officer, Naval Facilities Engineering Command, Hawaii 400 Marshall Road JBPHH, HI 96860-3139

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ACRONYMS AND ABBREVIATIONS

| ACRONYMS/ ABBREVIATIONS | DEFINITION/MEANING |
|----------------------------|---|
| % COPC DLNR | percent Contaminant of Potential Concern State of Hawaii Department of Land and Natural Resources |
| DOH | State of Hawaii Department of Land and Natural Resources State of Hawaii Department of Health |
| DON EAL | Department of the Navy Environmental Action Level |
| EPA | Environmental Protection Agency |
| ESI F-76 | Environmental Science International |
| ID | Marine Diesel Fuel Identification |
| JBPHH | Joint Base Pearl Harbor-Hickam |
| JP-5 JP-8 | Jet Fuel Propellant-5 Jet Fuel Propellant-8 |
| LCS | Laboratory Control Sample |
| LCSD | Laboratory Control Sample Duplicate |
| LOD LOQ | Limit of Detection Limit of Quantitation |
| µg/L | micrograms per liter |
| MS MSD | Matrix Spike Matrix Spike Duplicate |
| NAVFAC | Naval Facilities Engineering Command |
| NAVSUP FLC | Naval Supply Systems Command Fleet Logistics Center |
| N.D. PAH | Not Detected Polycyclic Aromatic Hydrocarbons |
| PARCCS | Precision, Accuracy, Representativeness, Completeness, Comparability, |
| рН | and Sensitivity hydrogen activity |
| QC | Quality Control |
| RHSF | Red Hill Bulk Fuel Storage Facility |
| RPD SAP | Relative Percent Difference Sampling and Analysis Plan |
| TEC | The Environmental Company, Inc. |
| TPH-d TPH-g | Total Petroleum Hydrocarbons as diesel Total Petroleum Hydrocarbons as gasoline |
| U.S. | United States of America |
| UST VOC | Underground Storage Tank |
| WP | Volatile Organic Compounds Work Plan |
| | |

EXECUTIVE SUMMARY

This quarterly monitoring report presents the results of the third quarter 2013 groundwater sampling event conducted on July 24, 2013, at the outside tunnel wells of the Red Hill Bulk Fuel Storage Facility [RHSF], Joint Base Pearl Harbor-Hickam [JBPHH], Hawaii. The RHSF is located in Halawa Heights on the Island of Oahu. There are 18 active and 2 inactive underground storage tanks [USTs] located at the RHSF. The State of Hawaii Department of Health [DOH] Facility Identification [ID] number is 9-102271. The DOH Release ID numbers are 990051, 010011, and 020028.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring at the RHSF, under Naval Facilities Engineering Command [NAVFAC] Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved Work Plan [WP]/Sampling and Analysis Plan [SAP] prepared by Environmental Science International [ESI].

On July 24, 2013, ESI personnel collected groundwater samples from two outside tunnel monitoring wells (wells HDMW2253-03 and OWDFMW01). A summary of the analytical results is provided below.

- **HDMW2253-03** Naphthalene (0.03 micrograms per liter [µg/L]) was detected. The Contaminant of Potential Concern [COPC] was not detected at a concentration above the DOH Environmental Action Levels [EALs] for drinking water toxicity or gross contamination.
- OWDFMW01 Total Petroleum Hydrocarbons as diesel [TPH-d] (470 and 340 μg/L), acetone (88 and 83 μg/L), naphthalene (0.081 and 0.12 μg/L), and benzene (0.42 and 0.44 μg/L) were detected in both the primary and duplicate samples. TPH-d (470 and 340 μg/L) was detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination in both samples.

TPH-d was not detected in well HDMW2253-03 during this round of quarterly sampling. TPH-d concentrations last exceeded the DOH EALs for both drinking water toxicity and gross contamination in January 2013 (600 μ g/L). Naphthalene was detected but remained at a low concentration, below the DOH EALs and consistent with historical results.

With the exception of TPH-d in well OWDFMW01, groundwater contaminant concentrations remained at low concentrations and did not change significantly, or were not detected. TPH-d concentrations detected during this round of sampling were above the DOH EALs for both drinking water toxicity and gross contamination, but decreased significantly from the last event, and were the lowest they have been since July 2012.

Based on the results of the assessment, continued groundwater monitoring at the RHSF is recommended. If the TPH-d concentrations significantly increase, the monitoring frequency

should be increased to monthly, even though the two outside wells are not included in the RHSF Groundwater Protection Plan.

SECTION 1 – INTRODUCTION

This quarterly monitoring report presents the results of the third quarter 2013 groundwater sampling event conducted on July 24, 2013, at the outside tunnel wells of the RHSF, JBPHH, Hawaii. The RHSF is located in Halawa Heights on the Island of Oahu. The purpose of the sampling is to (1) assess the condition of groundwater beneath and in the vicinity of the RHSF with respect to chemical constituents associated with jet fuel propellant and marine diesel fuel, and (2) to ensure the Navy remains in compliance with DOH UST release response requirements as described in Hawaii Administrative Rules 11-281 Subchapter 7, Release Response Action (DOH, 2000). The DOH Facility ID number for the RHSF is 9-102271. The DOH Release ID numbers are 990051, 010011, and 020028.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI (ESI, 2012).

1.1 SITE DESCRIPTION

The RHSF is located on federal government land (zoned F1- Military and Federal), located in Halawa Heights, approximately 2.5 miles northeast of Pearl Harbor (Figure 1). It is located on a low ridge on the western edge of the Koolau Mountain Range that divides Halawa Valley from Moanalua Valley. The RHSF is bordered on the north by Halawa Correctional Facility and private businesses, on the west by the United States of America [U.S.] Coast Guard reservation, on the south by residential neighborhoods, and on the east by Moanalua Valley. A quarry is located less than a quarter mile away to the northwest. The RHSF occupies 144 acres of land and the majority of the site is at an elevation of approximately 200 to 500 feet above mean sea level.

The RHSF contains 18 active and 2 inactive USTs, which are operated by Naval Supply Systems Command Fleet Logistics Center [NAVSUP FLC] Pearl Harbor (formerly Fleet and Industrial Supply Center). Each UST has a capacity of approximately 12.5 million gallons. The RHSF is located approximately 100 feet above the basal aquifer. The USTs contain Jet Fuel Propellant-5 [JP-5], Jet Fuel Propellant-8 [JP-8], and Marine Diesel Fuel [F-76]. The current status of each of the USTs is summarized in Table 1.1.

Two groundwater monitoring wells (well HDMW2253-03 and OWDFMW01) are located outside of the RHSF tunnel system (Figure 2). Well HDMW2253-03 is located at the Halawa Correctional Facility (outside the RHSF) and well OWDFMW01 is located at the Oily Waste Disposal Facility near Adit 3. Five groundwater monitoring wells (wells RHMW01, RHMW02, RHMW03, RHMW05, and RHMW2254-01) are located within the RHSF lower access tunnel. Monitoring data for the five wells located inside the tunnel are included in a separate report.

Monitoring wells RHMW01, RHMW02, RHMW03, and RHMW05 are located inside the underground tunnels. Monitoring well RHMW2254-01 is located inside the infiltration gallery of

the Department of the Navy [DON] Well 2254-01. DON Well 2254-01 is located approximately 2,400 feet downgradient of the USTs and provides approximately 24 percent [%] of the potable water to the Pearl Harbor System, which serves approximately 52,200 military customers. NAVFAC Public Works Department operates the infiltration gallery and DON Well 2254-01.

| Red Hill Bulk Fuel Storage Facility April 2013 Quarterly Monitoring Report | | | | | | | | | | | |
|--|-----------|----------|----------------------|--|--|--|--|--|--|--|--|
| Tank Identification | Fuel Type | Status | Capacity | | | | | | | | |
| F-1 | None | Inactive | 12.5 million gallons | | | | | | | | |
| F-2 | JP-8 | Active | 12.5 million gallons | | | | | | | | |
| F-3 | JP-8 | Active | 12.5 million gallons | | | | | | | | |
| F-4 | JP-8 | Active | 12.5 million gallons | | | | | | | | |
| F-5 | JP-8 | Active | 12.5 million gallons | | | | | | | | |
| F-6 | JP-8 | Active | 12.5 million gallons | | | | | | | | |
| F-7 | JP-5 | Active | 12.5 million gallons | | | | | | | | |
| F-8 | JP-5 | Active | 12.5 million gallons | | | | | | | | |
| F-9 | JP-5 | Active | 12.5 million gallons | | | | | | | | |
| F-10 | JP-5 | Active | 12.5 million gallons | | | | | | | | |
| F-11 | JP-5 | Active | 12.5 million gallons | | | | | | | | |
| F-12 | JP-5 | Active | 12.5 million gallons | | | | | | | | |
| F-13 | F-76 | Active | 12.5 million gallons | | | | | | | | |
| F-14 | F-76 | Active | 12.5 million gallons | | | | | | | | |
| F-15 | F-76 | Active | 12.5 million gallons | | | | | | | | |
| F-16 | F-76 | Active | 12.5 million gallons | | | | | | | | |
| F-17 | JP-5 | Active | 12.5 million gallons | | | | | | | | |
| F-18 | JP-5 | Active | 12.5 million gallons | | | | | | | | |
| F-19 | None | Inactive | 12.5 million gallons | | | | | | | | |
| F-20 | JP-5 | Active | 12.5 million gallons | | | | | | | | |

TABLE 1.1 Current Status of the USTs Red Hill Bulk Fuel Storage Facility April 2013 Quarterly Monitoring Report

F-76 Marine Diesel Fuel

JP-5 Jet Fuel Propellant-5

JP-8 Jet Fuel Propellant-8

1.2 PHYSICAL SETTINGS

Climatological conditions in the area of the RHSF consist of warm to moderate temperatures and low to moderate rainfall. The RHSF is leeward of the prevailing northeasterly trade winds. The average annual precipitation is approximately 40 inches, which occurs mainly between November and April (State of Hawaii Department of Land and Natural Resources [DLNR], 1986). Annual pan evaporation is approximately 75 inches (DLNR, 1985). Average temperatures range from the low 60's to high 80's (degrees Fahrenheit) (Atlas of Hawaii, 1983).

Oahu consists of the eroded remnants of two shield volcanoes, Waianae and Koolau. The RHSF is located on the southwest flank of the Koolau volcanic shield. Lavas erupted during the shield-building phase of the volcano belong to the *Koolau Volcanic Series* (Stearns and Vaksvik, 1935). Following formation of the Koolau shield, a long period of volcanic quiescence occurred, during which the shield was deeply eroded. Following this erosional period, eruptive activity resumed. Lavas and pyroclastic material erupted during this period belong to the *Honolulu Volcanic Series* (Stearns and Vaksvik, 1935).

In the immediate area of the RHSF, Koolau Volcanic Series lavas dominate, although there are consolidated and unconsolidated non-calcareous deposits in the vicinity that consist of alluvium generated during erosion of the Koolau volcanic shield. South-southwest of the Site, and in isolated exposures to the west, are pyroclastic deposits formed during eruptions from three Honolulu Volcanic Series vents, Salt Lake, Aliamanu, and Makalapa (Stearns and Vaksvik, 1935). Based on established geology and records of the drilled wells (Stearns and Vaksvik, 1938), the RHSF is underlain by Koolau Volcanic Series basalts. The area of the RHSF is classified as *Rock Land*, where 25 to 90% of the land surface is covered by exposed rock and there are only shallow soils (Foote, et al., 1972).

Groundwater in Hawaii exists in two principal types of aquifers. The first and most important type, in terms of drinking water resources, is the basal aquifer. The basal aquifer exists as a lens of fresh water floating on and displacing seawater within the pore spaces, fractures, and voids of the basalt that forms the underlying mass of each Hawaiian island. In parts of Oahu, groundwater in the basal aquifer is confined by the overlying caprock and is under pressure. Waters that flow freely to the surface from wells that tap the basal aquifer are referred to as *artesian*.

The second type of aquifer is the caprock aquifer, which consists of various kinds of unconfined and semi-confined groundwater. Commonly, the caprock consists of a thick sequence of nearly impermeable clays, coral, and basalt, which separates the caprock aquifer from the basal aquifer. The impermeable nature of these materials and the artesian nature of the basal aquifer severely restrict the downward migration of groundwater from the upper caprock aquifer. In the area of the RHSF, there is no discernible caprock.

Groundwater in the area of the RHSF is part of the *Waimalu Aquifer System* of the *Pearl Harbor Aquifer Sector*. The aquifer is classified as a basal, unconfined, flank-type; and is currently used as a drinking water source. The aquifer is considered fresh with less than 250 milligrams per liter of chloride and is considered an irreplaceable resource with a high vulnerability to contamination (Mink and Lau, 1990).

The nearest drinking water supply well is the DON Well 2254-01, located in the infiltration gallery within the RHSF. DON Well 2254-01 is located approximately 2,400 feet downgradient of the USTs (Figure 2).

1.3 BACKGROUND

The RHSF was constructed by the U.S. Government in the early 1940s. Twenty USTs and a series of tunnels were constructed to supply fuel to the Navy. The USTs were constructed of steel and they currently contain JP-5, JP-8, and F-76. Several tanks in the past have stored DON special fuel oil, DON distillate, aviation gasoline, and motor gasoline (Environet, 2010). The fueling system is a self-contained underground unit that was installed into native rock comprised primarily of basalt with some interbedded tuffs and breccias (Environet, 2010). Each

UST measures approximately 245 feet in height and 100 feet in diameter. The upper domes of the tanks lie at depths varying between 100 feet and 200 feet below ground surface.

In response to increasing concentrations of COPCs in the groundwater monitoring wells within the facility (specifically RHMW02) during the 2008 sampling events, quarterly groundwater monitoring was initiated in 2009 at the outside tunnel wells.

In 2009, groundwater samples were collected from wells RHMW04, OWDFMW01, and HDMW2253-03. Samples were collected in August and October 2009. None of the COPCs were detected at concentrations exceeding the gross contamination or drinking water toxicity DOH EALs.

In 2010, groundwater samples were collected from wells RHMW04, OWDFMW01, and HDMW2253-03. Samples were collected from well RHMW04 in January and April 2010. Samples were collected from well OWDFMW01 in January, April, and October 2010. Samples were collected from well HDMW2253-03 in January, April, July and October 2010. The COPCs concentrations exceeding DOH EALs are summarized below.

- **HDMW2253-03** TPH-d was detected at a concentration above the gross contamination and drinking water toxicity DOH EAL in January 2010 (The Environmental Company, Inc. [TEC], 2010a).
- **OWDFMW01** TPH-d was detected at a concentration above the gross contamination and drinking water toxicity DOH EALs in January and April 2010 (TEC, 2010a; TEC, 2010b).

In 2011, groundwater samples were collected from wells OWDFMW01 and HDMW2253-03. Samples were collected in January, April, July, and October 2011. None of the COPCs were detected at concentrations exceeding the gross contamination or drinking water toxicity DOH EALs.

In 2012, groundwater samples were collected from wells OWDFMW01 and HDMW2253-03. Samples were collected in January, April, July, and November 2012. TPH-d was detected at a concentration above the DOH EALs in samples collected from wells HDMW2253-03 and OWDFMW01 (Environet, 2012; ESI, 2013a). The COPCs concentrations exceeding DOH EALs are summarized below.

- **HDMW2253-03** TPH-d was detected at a concentration above the DOH EALs for gross contamination and drinking water toxicity in April and November 2012.
- **OWDFMW01** TPH-d was detected at a concentration above the DOH EALs for gross contamination and drinking water toxicity in April 2012.

In January 2013, groundwater samples were collected from wells OWDFMW01 and HDMW2253-03 (ESI 2013b). TPH-d was detected at a concentration above the DOH EALs in samples collected from wells HDMW2253-03 and OWDFMW01. The COPCs concentrations exceeding DOH EALs are summarized below.

- **HDMW2253-03** TPH-d was detected at a concentration above the DOH EALs for gross contamination and drinking water toxicity in January 2013.
- **OWDFMW01** TPH-d was detected at a concentration above the DOH EALs for gross contamination and drinking water toxicity in January 2013.

In April 2013, groundwater samples were collected from wells OWDFMW01 and HDMW2253-03 (ESI 2013c). TPH-d was detected at a concentration above the DOH EALs in samples collected from well OWDFMW01. The COPC concentrations exceeding DOH EALs are summarized below.

• **OWDFMW01** – TPH-d was detected at a concentration above the DOH EALs for gross contamination and drinking water toxicity in April 2013.

1.3.1 Previous Reports

The following groundwater monitoring reports were previously submitted to the DOH:

- 1. Groundwater Monitoring Report, August 2009 (submitted September 2009).
- 2. Groundwater Monitoring Report, October 2009 (submitted December 2009).
- 3. Groundwater Monitoring Report, January, 2010 (submitted April 2010).
- 4. Groundwater Monitoring Report, April 2010 (submitted May 2010).
- 5. Groundwater Monitoring Report, July 2010 (submitted August 2010).
- 6. Groundwater Monitoring Report, October 2010 (submitted December 2010).
- 7. Groundwater Monitoring Report, January 2011 (submitted March 2011).
- 8. Groundwater Monitoring Report, April 2011 (submitted June 2011).
- 9. Groundwater Monitoring Report, July 2011 (submitted September 2011).
- 10. Groundwater Monitoring Report, October 2011 (submitted December 2011).
- 11. Groundwater Monitoring Report, January 2012 (submitted March 2012).
- 12. Groundwater Monitoring Report, April 2012 (submitted July 2012).
- 13. Groundwater Monitoring Report, July 2012 (submitted August 2012).
- 14. Groundwater Monitoring Report, November 2012 (submitted January 2013).

- 15. Groundwater Monitoring Report, January 2013 (submitted April 2013).
- 16. Groundwater Monitoring Report, April 2013 (submitted July 2013).

SECTION 2 – GROUNDWATER SAMPLING

On July 24, 2013, ESI personnel collected groundwater samples from two monitoring wells (wells OWDFMW01 and HDMW2253-03). The samples were collected in accordance with DOH UST release response requirements and the RHSF Groundwater Protection Plan (TEC, 2008). Prior to purging and sampling, the depth to groundwater and the depth to the bottom of the wells were measured. Well OWDFMW01 was measured by ESI using a Geotech oil/water interface probe. Well HDMW2253-03 was measured by the DLNR using their Geotech oil/water interface probe. The measurements are included in the groundwater sampling logs. No measurable product, sheen, or petroleum hydrocarbon odor was observed in either well.

2.1 GROUNDWATER SAMPLING

Prior to collecting groundwater samples, disposable bailers were used to purge groundwater from the monitoring wells. Wells OWDFMW01 and HDMW2253-03 were purged at a rate of 0.29 and 0.11 liters per minute, respectively.

Water quality parameters were monitored on a periodic basis during well purging. The water quality parameters that were measured included hydrogen activity [pH], temperature, conductivity, dissolved oxygen, and oxidation reduction potential. The water quality parameters were evaluated to demonstrate that the natural characteristics of the aquifer formation water were present within the monitoring well before collecting the sample. At least four readings were collected during the purging process. Purging was considered complete when at least three consecutive water quality measurements stabilized within approximately 10%. The readings were recorded on groundwater monitoring logs which are included in Appendix A. The field notes are included in Appendix B.

When the water quality parameters stabilized, groundwater samples were collected from the wells. The disposable bailers were used to collect the groundwater samples from the monitoring wells. For each monitoring well, the groundwater samples were collected no more than two hours after purging was completed to prevent groundwater interaction with the monitoring well casing and atmosphere. Samples collected for dissolved lead were filtered in the field using a peristaltic pump and a 0.45 micron filter.

2.2 ANALYTICAL RESULTS

The samples were analyzed for TPH-d using U.S. Environmental Protection Agency [EPA] Method 8015M, Total Petroleum Hydrocarbons as gasoline [TPH-g] and Volatile Organic Compounds [VOCs] using EPA Method 8260B, Polycyclic Aromatic Hydrocarbons [PAHs] using EPA Method 8270C SIM, and dissolved lead using EPA Method 6020. The analytical results are summarized below and in Table 2.1. A copy of the laboratory report is included in Appendix C.

HDMW2253-03 – Naphthalene (0.03 μg/L) was detected. The COPC was not detected at a concentration above the DOH EALs.

OWDFMW01 – TPH-d (470 and 340 μg/L), acetone (88 and 83 μg/L), naphthalene (0.081 and 0.012 μg/L), and benzene (0.42 and 0.44 μg/L) were detected in both the primary and duplicate samples. TPH-d (470 and 340 μg/L) was detected at a concentration above the DOH EALs for both drinking water toxicity and gross contamination in both samples.

2.2.1 Groundwater Contaminant Trends

Historical groundwater contaminant concentration trends of COPCs that exceed the DOH EALs are presented in Appendix D. A summary of groundwater contaminant trends is provided below.

- HDMW2253-03 TPH-d was not detected during this round of quarterly sampling. TPH-d concentrations last exceeded the DOH EALs for both drinking water toxicity and gross contamination in January 2013 (600 µg/L). Naphthalene was detected but remained at a low concentration, below the DOH EALs and consistent with historical results.
- **OWDFMW01** With the exception of TPH-d, groundwater contaminant concentrations remained at low concentrations and did not change significantly, or were not detected. TPH-d concentrations detected during this round of sampling were above the DOH EALs for both drinking water toxicity and gross contamination, but decreased significantly from the last event, and were the lowest they have been since July 2012.

2.3 WASTE DISPOSAL

The purged groundwater and decontamination water generated during sampling of the wells was stored in a 55-gallon drum along with the purged water and decontamination water from the inside tunnel wells. The drum was stored onsite at Adit 3. On August 23, 2013, the drum of water was picked up by Pacific Commercial Services, LLC and disposed at Unitek Solvent Services, Inc. The waste disposal manifest is included in Appendix E.

TABLE 2.1 Analytical Results for Groundwater Sampling (July 24, 2013) Red Hill Bulk Fuel Storage Facility July 2013 Quarterly Monitoring Report

| EPA 8260B T A | Chemical TPH-d TPH-g | DOH Drinking Water Toxicity | EALs Gross | | OW | DFMW01 (ES | J34) | | | OWDF | MW01 (ES035 | b) (Dup) | | | HDM | W2253-03 (ES | 5036) | |
|--------------------------------------|---|-----------------------------------|---------------|---------------|-----|------------------|------------|-------------|--------------|------|-------------|-------------|-------------|--------------|----------|--------------|-------------|----------------|
| EPA 8015B T EPA 8260B T A A | TPH-d | | Gross | | | OWDFMW01 (ES034) | | | | | | | | | | | | |
| EPA 8260B T A | | | Contamination | Results | Q | LOQ | LOD | DL | Results | Q | LOQ | LOD | DL | Results | Q | LOQ | LOD | DL |
| A A | IPH-g | 190 | 100 | 470 | HD | 52 | 21 | 15 | 340 | HD | 52 | 21 | 15 | N.D. | U | 52 | 21 | 15 |
| A | Assasshthana | 100 | 100 | N.D. N.D. | UU | 50 | 30 0.05 | 13 0.021 | N.D. N.D. | U | 50 0.19 | 30 0.048 | 13 0.020 | N.D. N.D. | U U | 50 0.19 | 30 0.047 | 13 0.020 |
| | Acenaphthene Acenaphthylene | 370 240 | 20 2,000 | N.D. N.D. | U | 0.2 | 0.05 | 0.021 | N.D. N.D. | UU | 0.19 | 0.048 | 0.020 | N.D. N.D. | U | 0.19 | 0.047 | 0.020 |
| | Anthracene | 1.800 | 2,000 | N.D. | U | 0.2 | 0.05 | 0.018 | N.D. | U | 0.19 | 0.048 | 0.033 | N.D. | <u> </u> | 0.19 | 0.047 | 0.032 |
| | Benzo[a]anthracene | 0.092 | 4.7 | N.D. | U | 0.2 | 0.05 | 0.024 | N.D. | U | 0.19 | 0.048 | 0.023 | N.D. | <u> </u> | 0.19 | 0.047 | 0.022 |
| | Benzo[g,h,i]perylene | 1,500 | 0.13 | N.D. | U | 0.2 | 0.05 | 0.022 | N.D. | U | 0.19 | 0.048 | 0.020 | N.D. | U | 0.19 | 0.047 | 0.021 |
| | Benzo[a]pyrene | 0.2 | 0.81 | N.D. | U | 0.2 | 0.05 | 0.036 | N.D. | U | 0.19 | 0.048 | 0.035 | N.D. | U | 0.19 | 0.047 | 0.034 |
| | Benzo[b]fluoranthene | 0.092 | 0.75 | N.D. | U | 0.2 | 0.05 | 0.025 | N.D. | U | 0.19 | 0.048 | 0.024 | N.D. | U | 0.19 | 0.047 | 0.024 |
| E | Benzo[k]fluoranthene | 0.92 | 0.4 | N.D. | U | 0.2 | 0.05 | 0.023 | N.D. | U | 0.19 | 0.048 | 0.023 | N.D. | U | 0.19 | 0.047 | 0.022 |
| EPA 8270C SIM | Chrysene | 9.2 | 1 | N.D. | U | 0.2 | 0.05 | 0.019 | N.D. | U | 0.19 | 0.048 | 0.018 | N.D. | U | 0.19 | 0.047 | 0.018 |
| L | Dibenzo[a,h]anthracene | 0.0092 | 0.52 | N.D. | U | 0.2 | 0.05 | 0.027 | N.D. | U | 0.19 | 0.048 | 0.026 | N.D. | U | 0.19 | 0.047 | 0.025 |
| | Fluoranthene | 1,500 | 130 | N.D. | U | 0.2 | 0.05 | 0.027 | N.D. | U | 0.19 | 0.048 | 0.026 | N.D. | U | 0.19 | 0.047 | 0.026 |
| | Fluorene | 240 | 950 | N.D. | U | 0.2 | 0.05 | 0.024 | N.D. | U | 0.19 | 0.048 | 0.024 | N.D. | U | 0.19 | 0.047 | 0.023 |
| | Indeno[1,2,3-cd]pyrene | 0.092 | 0.095 | N.D. | U | 0.2 | 0.05 | 0.022 | N.D. | U | 0.19 | 0.048 | 0.021 | N.D. | U | 0.19 | 0.047 | 0.021 |
| | 1,-Methylnaphthalene | 4.7 | 10 10 | N.D. N.D. | UU | 0.2 | 0.05 | 0.028 | N.D. N.D. | UU | 0.19 | 0.048 | 0.027 | N.D. N.D. | U U | 0.19 0.19 | 0.047 | 0.027 0.025 |
| | 2,-Methylnaphthalene Naphthalene | 17 | 21 | N.D. 0.081 | J | 0.2 | 0.05 | 0.026 | 0.12 | J | 0.19 | 0.048 | 0.026 | 0.030 | | 0.19 | 0.047 | 0.025 |
| | Phenanthrene | 240 | 410 | N.D. | U | 0.2 | 0.05 | 0.030 | N.D. | Ű | 0.19 | 0.048 | 0.022 | N.D. | Ű | 0.19 | 0.047 | 0.029 |
| | Pyrene | 180 | 68 | N.D. | U U | 0.2 | 0.05 | 0.025 | N.D. | U | 0.19 | 0.048 | 0.023 | N.D. | <u> </u> | 0.19 | 0.047 | 0.023 |
| | 1,1,1-Trichloroethane | 200 | 970 | N.D. | U | 5.0 | 0.5 | 0.30 | N.D. | U | 5.0 | 0.5 | 0.30 | N.D. | U | 5.0 | 0.5 | 0.30 |
| | 1,1,2-Trichloroethane | 5 | 50,000 | N.D. | U | 1.0 | 0.5 | 0.38 | N.D. | U | 1.0 | 0.5 | 0.38 | N.D. | U | 1.0 | 0.5 | 0.38 |
| 1 | 1,1-Dichloroethane | 2.4 | 50,000 | N.D. | U | 5.0 | 0.5 | 0.28 | N.D. | U | 5.0 | 0.5 | 0.28 | N.D. | U | 5.0 | 0.5 | 0.28 |
| | 1,1-Dichloroethylene | 7 | 1,500 | N.D. | U | 1.0 | 0.5 | 0.43 | N.D. | U | 1.0 | 0.5 | 0.43 | N.D. | U | 1.0 | 0.5 | 0.43 |
| | 1,2,3-Trichloropropane | 0.6 | 50,000 | N.D. | U | 5.0 | 1.0 | 0.64 | N.D. | U | 5.0 | 1.0 | 0.64 | N.D. | U | 5.0 | 1.0 | 0.64 |
| | 1,2,4-Trichlorobenzene | 70 | 3,000 | N.D. | U | 5.0 | 1.0 | 0.5 | N.D. | U | 5.0 | 1.0 | 0.5 | N.D. | U | 5.0 | 1.0 | 0.5 |
| | 1,2-Dibromo-3- chloropropane | 0.04 | 10 | N.D. | U | 10 | 2.0 | 1.2 | N.D. | U | 10 | 2.0 | 1.2 | N.D. | U | 10 | 2.0 | 1.2 |
| | 1,2-Dibromoethane | 0.04 | 50,000 | N.D. | U | 1.0 | 0.5 | 0.36 | N.D. | U | 1.0 | 0.5 | 0.36 | N.D. N.D. | U | 1.0 | 0.5 | 0.36 |
| | 1,2-Dichlorobenzene 1,2-Dichloroethane | 600 0.15 | 10 7,000 | N.D. N.D. | UU | 1.0 | 0.5 0.5 | 0.46 | N.D. N.D. | UU | 1.0 | 0.5 | 0.46 | N.D. N.D. | U U | 1.0 1.0 | 0.5 | 0.46 |
| | 1,2-Dichloropropane | 5 | 10 | N.D. | U | 5.0 | 0.5 | 0.42 | N.D. | U | 5.0 | 0.5 | 0.24 | N.D. | U | 5.0 | 0.5 | 0.42 |
| | 1,3-Dichlorobenzene | 180 | 5 | N.D. | U | 1.0 | 0.5 | 0.4 | N.D. | Ŭ | 1.0 | 0.5 | 0.4 | N.D. | U | 1.0 | 0.5 | 0.4 |
| | 1,3-Dichloropropene (total of cis/trans) | 0.43 | 50,000 | N.D. | U | 1.0 | 0.5 | 0.25 | N.D. | U | 1.0 | 0.5 | 0.25 | N.D. | U | 1.0 | 0.5 | 0.25 |
| 1 | 1,4-Dichlorobenzene | 75 | 5 | N.D. | U | 1.0 | 0.5 | 0.43 | N.D. | U | 1.0 | 0.5 | 0.43 | N.D. | U | 1.0 | 0.5 | 0.43 |
| A | Acetone | 22,000 | 20,000 | 88 | | 20 | 10 | 6.0 | 83 | | 20 | 10 | 6.0 | N.D. | U | 20 | 10 | 6.0 |
| E | Benzene | 5 | 170 | 0.42 | J | 1.0 | 0.5 | 0.14 | 0.44 | J | 1.0 | 0.5 | 0.14 | N.D. | U | 1.0 | 0.5 | 0.14 |
| | Bromodichloromethane | 0.12 | 50,000 | N.D. | U | 5.0 | 0.5 | 0.21 | N.D. | U | 5.0 | 0.5 | 0.21 | N.D. | U | 5.0 | 0.5 | 0.21 |
| | Bromoform | 80 | 510 | N.D. | U | 10 | 1.0 | 0.50 | N.D. | U | 10 | 1.0 | 0.50 | N.D. | U | 10 | 1.0 | 0.50 |
| | Bromomethane | 8.7 | 50,000 | N.D. | U | 20 | 5.0 | 3.9 | N.D. | U | 20 | 5.0 | 3.9 | N.D. | U | 20 | 5.0 | 3.9 |
| | Carbon Tetrachloride | 5 100 | 520 | N.D. N.D. | U | 1.0 | 0.5 | 0.23 | N.D. | U | 1.0 | 0.5 | 0.23 | N.D. | U | 1.0 | 0.5 | 0.23 |
| | Chlorobenzene Chloroethane | 21,000 | 50 16 | N.D. | UU | 5.0 10 | 0.5 5.0 | 0.17 2.3 | N.D. N.D. | UU | 5.0 10 | 5.0 | 0.17 2.3 | N.D. N.D. | U U | 5.0 10 | 0.5 5.0 | 0.17 2.3 |
| | Chloroform | 70 | 2,400 | N.D. | U | 5.0 | 0.5 | 0.46 | N.D. | U | 5.0 | 0.5 | 0.46 | N.D. | U | 5.0 | 0.5 | 0.46 |
| | Chloromethane | 1.8 | 50,000 | N.D. | U U | 10 | 2.0 | 1.8 | N.D. | U | 10 | 2.0 | 1.8 | N.D. | <u> </u> | 10 | 2.0 | 1.8 |
| | cis-1,2-Dichloroethylene | 70 | 50,000 | N.D. | U | 1.0 | 0.5 | 0.48 | N.D. | U | 1.0 | 0.5 | 0.48 | N.D. | U | 1.0 | 0.5 | 0.48 |
| | Dibromochloromethane | 0.16 | 50,000 | N.D. | U | 1.0 | 0.5 | 0.25 | N.D. | U | 1.0 | 0.5 | 0.25 | N.D. | U | 1.0 | 0.5 | 0.25 |
| E | Ethylbenzene | 700 | 30 | N.D. | U | 1.0 | 0.5 | 0.14 | N.D. | U | 1.0 | 0.5 | 0.14 | N.D. | U | 1.0 | 0.5 | 0.14 |
| | Hexachlorobutadiene | 0.86 | 6 | N.D. | U | 1.0 | 0.5 | 0.32 | N.D. | U | 1.0 | 0.5 | 0.32 | N.D. | U | 1.0 | 0.5 | 0.32 |
| | Methyl ethyl ketone (2-Butanone) | 7,100 | 8,400 | N.D. | U | 10 | 5.0 | 2.2 | N.D. | U | 10 | 5.0 | 2.2 | N.D. | U | 10 | 5.0 | 2.2 |
| | Methyl isobutyl ketone (4-Methyl-2-Pentanone) | 2,000 | 1300 | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 |
| | Methyl tert-butyl Ether | 12 4.8 | 5 9,100 | N.D. N.D. | U | 1.0 | 0.5 | 0.31 | N.D. | UU | 1.0 | 0.5 | 0.31 | N.D. N.D. | U | 1.0 | 0.5 | 0.31 |
| | Methylene chloride Styrene | 4.8 | 9,100 | N.D. N.D. | UU | 5.0 1.0 | 1.0 0.5 | 0.64 | N.D. N.D. | U | 5.0 | 1.0 0.5 | 0.64 | N.D. N.D. | U U | 5.0 1.0 | 1.0 0.5 | 0.64 |
| | Tetrachloroethane, 1,1,1,2- | 0.52 | 50,000 | N.D. | U | 1.0 | 0.5 | 0.17 | N.D. | U | 1.0 | 0.5 | 0.17 | N.D. | U | 1.0 | 0.5 | 0.17 |
| | Tetrachloroethane, 1,1,2,2- | 0.067 | 500 | N.D. | UJ | 1.0 | 0.5 | 0.40 | N.D. | UJ | 1.0 | 0.5 | 0.40 | N.D. | UJ | 1.0 | 0.5 | 0.40 |
| | Tetrachloroethylene | 5 | 170 | N.D. | U | 5.0 | 0.5 | 0.39 | N.D. | U | 5.0 | 0.5 | 0.39 | N.D. | U | 5.0 | 0.5 | 0.39 |
| | Toluene | 1,000 | 40 | N.D. | U | 1.0 | 0.5 | 0.24 | N.D. | U | 1.0 | 0.5 | 0.24 | N.D. | U | 1.0 | 0.5 | 0.24 |
| ti | trans-1,2- Dichloroethylene | 100 | 260 | N.D. | U | 1.0 | 0.5 | 0.37 | N.D. | U | 1.0 | 0.5 | 0.37 | N.D. | U | 1.0 | 0.5 | 0.37 |
| Т | Trichloroethylene | 5 | 310 | N.D. | U | 1.0 | 0.5 | 0.37 | N.D. | U | 1.0 | 0.5 | 0.37 | N.D. | U | 1.0 | 0.5 | 0.37 |
| | Vinyl chloride | 2 | 3,400 | N.D. | U | 1.0 | 0.5 | 0.30 | N.D. | U | 1.0 | 0.5 | 0.30 | N.D. | U | 1.0 | 0.5 | 0.30 |
| | Xylenes | 10,000 | 20 | N.D. | U | 11 | 1.5 | 0.23 | N.D. | U | 11 | 1.5 | 0.23 | N.D. | U | 11 | 1.5 | 0.23 |
| EPA 6020 E | Dissolved Lead | 15 | 50,000 | N.D. | U | 1.0 | 0.2 | 0.0898 | N.D. | U | 1 | 0.2 | 0.0898 | N.D. | U | 1 | 0.2 | 0.0898 |

 Image: Disponse Lead
 15
 50,000
 N.D.
 U
 1.0
 0.2

 The data are in micrograms per liter (µg/L). Shaded values exceeded the DOH EALs.
 DOH EALs
 DOH Tier 1 Environmental Action Levels for groundwater where groundwater is a current drinking water source and surface water is greater than 150 meters from the site (DOH, Fall 2011).
 DL
 Detection Limit or Method Detection Limit (MDL)

 EPA
 Environmental Protection Agency
 HD
 The chromatographic pattern was inconsistent with the profile of the reference fuel standard.
 J
 Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated.
 LOD
 Limit of Detection

Limit of Quantitation Not Detected Qualifiers

LOQ N.D.

Q TPH-g TPH-d

Total Petroleum Hydrocarbons as gasoline Total Petroleum Hydrocarbons as diesel Undetected at DL and is reported as less than the LOD.

U

SECTION 3 – DATA QUALITY ASSESSMENT

A data quality assessment, which consists of a review of the overall groundwater sample collection and analysis process, was performed in order to determine whether the analytical data generated meet the quality objectives for the project. The field Quality Control [QC] program consisted of standardized sample collection and management procedures, and the collection of field duplicate samples, matrix spike samples, and trip blank samples. The laboratory quality assurance program consisted of the use of standard analytical methods and the preparation and analyses of Matrix Spike [MS]/Matrix Spike Duplicate [MSD] samples, surrogate spikes, blanks, Laboratory Control Samples [LCS] and Laboratory Control Sample Duplicate [LCSD].

3.1 Data Validation and Assessment

The objective of data validation is to provide data of known quality for project decisions. Data quality is judged in terms of Precision, Accuracy, Representativeness, Completeness, Comparability, and Sensitivity [PARCCS]. A number of factors may affect the quality of data, including: sample collection methods, sample analysis methods, and adherence to established procedures for sample collection, preservation, management, shipment, and analysis.

Precision

Precision is defined as the reproducibility of replicate measurements. Precision is evaluated by Relative Percentage Difference [RPD] of field duplicates and laboratory LCS/LCSD or MS/MSD results. Field duplicate and MS/MSD samples were collected at a rate of approximately 10% of project samples. Field duplicates were sent to the laboratory along with the primary samples.

The RPDs of detected analytes for the primary and field duplicate samples (ES034 and ES035) are provided in Table 3.1. A precision of less than 50% for duplicate pairs is required by the DoN Project Procedures Manual to be considered acceptable (DoN 2007). For this monitoring event, the RPDs for duplicate sample pairs were all within the acceptable range except for TPH-d (32%) and naphthalene (39%). In samples ES034 and ES035, TPH-d was detected at concentrations approximately double the DOH EALs, and naphthalene was detected approximately two order of magnitude below the DOH EALs. Therefore, this slight reduction in precision is unlikely to affect data usability. The naphthalene concentrations detected in the samples were below the respective limits of quantitation [LOQs] implying a higher uncertainty for these results than for values detected above the LOQs (i.e. estimated, J-flagged). Consequently, the assigned RPDs signified the anticipated decrease in precision is considered acceptable based on the sample duplicated evaluation.

The RPD of the MS/MSD results for benzo[b]fluoranthene, dibenzo[a,h]anthracene, flourene, phenanthrene, and 1,1,2,2-tetrachloroethane were above the acceptable maximum of 20%. Only 1,1,2,2-tetrachloroethane was above 50%, an RPD recommended in the NAVFAC Project

Procedures Manual [DON 2007]. None of these COPCs were detected during this sampling event or have been historically detected. Therefore, this is unlikely to affect data usability.

Accuracy

Accuracy is defined as the degree of conformity of a measurement to a standard or true value. Accuracy is evaluated through measurement of the percent recovery of an analyte in a reference standard or spiked sample. Accuracy limits for surrogates, laboratory control spike, MS, and MSD samples are established by the individual laboratory. The acceptance criteria for accuracy are dependent on the analytical method and are based on historical laboratory data.

Results for TPH-d in samples ES034 and ES035 were flagged "HD." The laboratory indicated a mismatch between the calibration standard and the TPH-d chromatographic profile. Mismatches of this type are not uncommon. The chromatograms are not part of the standard laboratory package and were not reviewed by ESI.

With the exception of acetone, all of the LCS and surrogate spike recoveries for analyzed constituents were within acceptable percent recovery limits. The MS and/or MSD recoveries were below the control limits for 1,1,2,2-tetrachloroethane and the associated sample results may be biased low. Sample results for 1,1,2,2-tetrachloroethane were flagged "UJ." 1,1,2,2-Tetrachloroethane was not detected in any of the groundwater samples. However the drinking water EAL for 1,1,2,2-Tetrachloroethane is below the respective limits of detection [LODs]. Based on historical results and the results of other VOCs in the sample; it is not likely that 1,1,2,2-tetrachloroethane is present at concentrations above the DOH EALs.

The MS and/or MSD recoveries were above the control limits for acetone, tetrachloroethylene, and trichloroethylene, and the associated sample results may be biased high. None of the COPCs were detected at concentrations above the DOH EALs, so a potential high bias should not affect data usability. The LCS recovery for acetone was also above the control limits.

All other MS/MSD recoveries were within acceptable recovery limits; therefore, the data accuracy for this monitoring event is considered acceptable.

Representativeness

Representativeness is the degree that data accurately and precisely represents a characteristic of a population, parameter variations at a sampling point, or an environmental condition. Representativeness was achieved by conducting sampling in compliance with the sample collection procedures described in the WP/SAP specifically written for this project (ESI, 2012).

Representativeness is also evaluated through the compliance with the sample holding time, sample preservation, and the analysis of blank samples, including method blank and trip blank samples. The sample holding time and sample preservation complied with the EPA criteria. For this sampling event, one trip blank was included in the cooler to assess contamination during sample transport for TPH-g and VOCs. TPH-g and VOCs were not detected in the trip blank.

Therefore, the groundwater sample data are considered representative of the groundwater quality on site.

Completeness

Completeness is defined as the overall percentage of valid analytical results (including estimated results) compared to the total number of analytical results reported by the laboratory. No data were rejected for this project, and therefore the completeness goal for this project (90%) was successfully met.

Comparability

Comparability expresses the confidence with which one data set can be compared to another data set. Comparability can be related to accuracy and precision because these quantities are measures of data reliability. Data, with acceptable precision and accuracy, are considered comparable if collection techniques, analytical procedures, methods and reporting are equivalent. For this monitoring event, the samples were collected using approaches consistent with those in the previous events, and the same analytical methods/procedures were used to measure the concentration of COPCs. Therefore, the results are considered comparable within this data set and with the data collected from previous sampling events. The field and laboratory personnel followed standard operation procedures.

All TPH-g data through July 2010 was analyzed by EPA Method 8015; beginning in October 2010, EPA Method 8260B was used. There was no event where both methods were used, and so there is no way to directly compare the results using each method and determine if one method produces biased results. However, there is no reason to believe that using either method should bias the data, and the TPH-g data for all events should be comparable.

Between August 2009 and July 2010, naphthalene was analyzed for by both EPA Methods 8260B and 8270C, and both results were reported. Beginning in October 2010, only results using EPA Method 8270C were reported. Naphthalene was not detected in any well until November 2012, so comparability with older results should not be a concern. However, the low bias associated with EPA Method 8270C must be considered when making project decisions.

Sensitivity

The LOQs are established by the laboratory based on the LODs or instrument detection limits, historical data, and EPA limits established for the methods. The LOQs for samples may require adjustment due to matrix interference or if high levels of target analytes necessitate dilution before analysis. Matrix interference and sample dilutions have the effect of increasing the LOQs. Laboratory LODs and LOQs for several analytes differed from the LODs and LOQs in the WP/SAP because the laboratory updates them quarterly. LODs and LOQs for several analytes were greater than the DOH EALs (as stated in the WP/SAP) and therefore it is not possible to determine whether the analytes are present at concentrations greater than or equal to the DOH EALs. As suggested by the DOH Technical Guidance Manual, the project action level will be the LOQ for these analytes.

3.2 Data Assessment and Usability Conclusions

The PARCCS criteria were evaluated, and with a few exceptions, all criteria were met. These exceptions include the exceedances of recovery criteria for MS/MSDs for several VOCs and PAHs. Since the surrogate recoveries and the recoveries of the VOCs and PAHs in the LCS/LCSD are all within recovery criteria, the MS/MSD exceendaces are not considered to affect the usability of the data, but may indicate some matrix heterogeneity. The data assessment concludes that all data generated during this event are usable for their intended purpose.

TABLE 3.1 Quality Control Results for Groundwater Sampling (July 24, 2013) Red Hill Bulk Fuel Storage Facility July 2013 Quarterly Monitoring Report

| | Chemical Constituent | DOH | EALs | | 0 | WDFMW01 (ES | | <u></u> | | OWDFMW01 (ES035) (DUP) | | | | | ES Trip | | | | |
|---------------|---|----------------|----------------------|--------------|----------|-------------|------|---------|--------------|------------------------|------------|-------|-----------|---------------------|--------------|---------|------------|---------|-----------|
| Method | | Drinking Water | Gross | Results | 0 | LOQ | LOD | DL | Results | Q | LOQ | LOD | DL | Duplicate | Results | Q | LOQ | LOD | DL |
| EPA 8015B | TPH-d | 190 | Contamination 100 | 470 | HD | 52 | 21 | 15 | 340 | HD | 52 | 21 | 15 | (%) 32.10 | - | | - | - | - |
| EPA 8260B | TPH-g | 100 | 100 | N.D. | U | 50 | 30 | 13 | N.D. | U | 50 | 30 | 13 | NA | N.D. | U | 50 | 30 | 13 |
| | Acenaphthene | 370 | 20 | N.D. | U | 0.2 | 0.05 | 0.021 | N.D. | U | 0.19 | 0.048 | 0.020 | NA | - | - | - | - | - |
| | Acenaphthylene | 240 | 2,000 | N.D. | U | 0.2 | 0.05 | 0.018 | N.D. | U | 0.19 | 0.048 | 0.017 | NA | - | - | - | - | - |
| | Anthracene | 1,800 | 22 | N.D. | U | 0.2 | 0.05 | 0.034 | N.D. | U | 0.19 | 0.048 | 0.033 | NA | - | - | - | - | - |
| | Benzo[a]anthracene | 0.092 | 4.7 | N.D. | U | 0.2 | 0.05 | 0.024 | N.D. | U | 0.19 | 0.048 | 0.023 | NA | - | - | - | - | - |
| | Benzo[g,h,i]perylene | 1,500 | 0.13 | N.D. | U | 0.2 | 0.05 | 0.022 | N.D. | U | 0.19 | 0.048 | 0.021 | NA | - | - | - | - | - |
| | Benzo[a]pyrene | 0.2 | 0.81 | N.D. N.D. | UU | 0.2 | 0.05 | 0.036 | N.D. N.D. | UU | 0.19 | 0.048 | 0.035 | NA NA | - | - | - | - | - |
| | Benzo[b]fluoranthene Benzo[k]fluoranthene | 0.92 | 0.4 | N.D. | U | 0.2 | 0.05 | 0.023 | N.D. | U | 0.19 | 0.048 | 0.024 | NA | - | - | - | - | - |
| | Chrysene | 9.2 | 1 | N.D. | U | 0.2 | 0.05 | 0.023 | N.D. | U | 0.19 | 0.048 | 0.025 | NA | - | - | | - | - |
| EPA 8270C SIM | Dibenzo[a,h]anthracene | 0.0092 | 0.52 | N.D. | U | 0.2 | 0.05 | 0.027 | N.D. | U | 0.19 | 0.048 | 0.026 | NA | - | - | - | - | - |
| | Fluoranthene | 1,500 | 130 | N.D. | U | 0.2 | 0.05 | 0.027 | N.D. | U | 0.19 | 0.048 | 0.026 | NA | - | - | - | - | - |
| | Fluorene | 240 | 950 | N.D. | U | 0.2 | 0.05 | 0.024 | N.D. | U | 0.19 | 0.048 | 0.024 | NA | - | - | - | - | - |
| 1 | Indeno[1,2,3-cd]pyrene | 0.092 | 0.095 | N.D. | U | 0.2 | 0.05 | 0.022 | N.D. | U | 0.19 | 0.048 | 0.021 | NA | - | - | - | - | - |
| 1 | 1,-Methylnaphthalene | 4.7 | 10 | N.D. | U | 0.2 | 0.05 | 0.028 | N.D. | U | 0.19 | 0.048 | 0.027 | NA | - | - | - | - | - |
| 1 | 2,-Methylnaphthalene | 24 | 10 | N.D. | U | 0.2 | 0.05 | 0.026 | N.D. | U | 0.19 | 0.048 | 0.026 | NA | - | - | - | - | - |
| 1 | Naphthalene | 17 | 21 | 0.081 | J | 0.2 | 0.05 | 0.023 | 0.12 | J | 0.19 | 0.048 | 0.022 | 38.81 | - | - | - | - | - |
| 1 | Phenanthrene Pyrene | 240 180 | 410 68 | N.D. N.D. | UU | 0.2 | 0.05 | 0.030 | N.D. N.D. | UU | 0.19 | 0.048 | 0.029 | NA NA | - | - | - | - | - |
| | 1,1,1-Trichloroethane | 200 | 970 | N.D. N.D. | U | 5.0 | 0.05 | 0.025 | N.D. | U | 5.0 | 0.048 | 0.024 | NA | - N.D. | - U | - 5.0 | - 0.5 | - 0.30 |
| | 1.1.2-Trichloroethane | 5 | 50,000 | N.D. | U U | 1.0 | 0.5 | 0.38 | N.D. | U | 1.0 | 0.5 | 0.38 | NA | N.D. | U | 1.0 | 0.5 | 0.38 |
| | 1.1-Dichloroethane | 2.4 | 50,000 | N.D. | U | 5.0 | 0.5 | 0.28 | N.D. | U | 5.0 | 0.5 | 0.28 | NA | N.D. | U | 5.0 | 0.5 | 0.28 |
| | 1,1-Dichloroethylene | 7 | 1,500 | N.D. | U | 1.0 | 0.5 | 0.43 | N.D. | U | 1.0 | 0.5 | 0.43 | NA | N.D. | U | 1.0 | 0.5 | 0.43 |
| | 1,2,3-Trichloropropane | 0.6 | 50,000 | N.D. | U | 5.0 | 1.0 | 0.64 | N.D. | U | 5.0 | 1.0 | 0.64 | NA | N.D. | U | 5.0 | 1.0 | 0.64 |
| | 1,2,4-Trichlorobenzene | 70 | 3,000 | N.D. | U | 5.0 | 1.0 | 0.5 | N.D. | U | 5.0 | 1.0 | 0.5 | NA | N.D. | U | 5.0 | 1.0 | 0.5 |
| | 1,2-Dibromo-3- chloropropane | 0.04 | 10 | N.D. | U | 10 | 2.0 | 1.2 | N.D. | U | 10 | 2.0 | 1.2 | NA | N.D. | U | 10 | 2.0 | 1.2 |
| | 1,2-Dibromoethane | 0.04 | 50,000 | N.D. | U | 1.0 | 0.5 | 0.36 | N.D. | U | 1.0 | 0.5 | 0.36 | NA | N.D. | U | 1.0 | 0.5 | 0.36 |
| | 1,2-Dichlorobenzene | 600 | 10 | N.D. | U | 1.0 | 0.5 | 0.46 | N.D. | U | 1.0 | 0.5 | 0.46 | NA | N.D. | U | 1.0 | 0.5 | 0.46 |
| | 1,2-Dichloroethane | 0.15 | 7,000 | N.D. | U | 1.0 | 0.5 | 0.24 | N.D. | U | 1.0 | 0.5 | 0.24 | NA | N.D. | U | 1.0 | 0.5 | 0.24 |
| | 1,2-Dichloropropane | 5 | 10 5 | N.D. | U | 5.0 | 0.5 | 0.42 | N.D. | U | 5.0 | 0.5 | 0.42 | NA | N.D. | U | 5.0 | 0.5 | 0.42 |
| | 1,3-Dichlorobenzene 1,3-Dichloropropene (total of cis/trans) | 180 | 50,000 | N.D. N.D. | UU | 1.0 | 0.5 | 0.4 | N.D. N.D. | UU | 1.0 | 0.5 | 0.4 0.25 | NA NA | N.D. N.D. | UU | 1.0 2.0 | 0.5 | 0.4 0.25 |
| | 1,4-Dichlorobenzene | 75 | 5 | N.D. | U U | 1.0 | 0.5 | 0.43 | N.D. | U | 1.0 | 0.5 | 0.43 | NA | N.D. | U | 1.0 | 0.5 | 0.23 |
| | Acetone | 22,000 | 20.000 | 88 | <u> </u> | 20 | 10 | 6.0 | 83 | <u> </u> | 20 | 10 | 6.0 | 5.85 | N.D. | U | 20 | 10 | 6.0 |
| | Benzene | 5 | 170 | 0.42 | J | 1.0 | 0.5 | 0.14 | 0.44 | J | 1.0 | 0.5 | 0.14 | 4.65 | N.D. | U | 1.0 | 0.5 | 0.14 |
| | Bromodichloromethane | 0.12 | 50,000 | N.D. | U | 5.0 | 0.5 | 0.21 | N.D. | U | 5.0 | 0.5 | 0.21 | NA | N.D. | U | 5.0 | 0.5 | 0.21 |
| | Bromoform | 80 | 510 | N.D. | U | 10 | 1.0 | 0.50 | N.D. | U | 10 | 1.0 | 0.50 | NA | N.D. | U | 10 | 1.0 | 0.50 |
| | Bromomethane | 8.7 | 50,000 | N.D. | U | 20 | 5.0 | 3.9 | N.D. | U | 20 | 5.0 | 3.9 | NA | N.D. | U | 20 | 5.0 | 3.9 |
| | Carbon Tetrachloride | 5 | 520 | N.D. | U | 1.0 | 0.5 | 0.23 | N.D. | U | 1.0 | 0.5 | 0.23 | NA | N.D. | U | 1.0 | 0.5 | 0.23 |
| EPA 8260B | Chlorobenzene | 100 | 50 | N.D. | U | 5.0 | 0.5 | 0.17 | N.D. | U | 5.0 | 0.5 | 0.17 | NA | N.D. | U | 5.0 | 0.5 | 0.17 |
| | Chloroethane | 21,000 | 16 | N.D. | U | 10 | 5.0 | 2.3 | N.D. | U | 10 | 5.0 | 2.3 | NA | N.D. | U | 10 | 5.0 | 2.3 |
| 1 | Chloroform Chloromethane | 70 | 2,400 50,000 | N.D. N.D. | UU | 5.0 | 0.5 | 0.46 | N.D. N.D. | UU | 5.0 10 | 0.5 | 0.46 | NA NA | N.D. N.D. | UU | 5.0 10 | 0.5 2.0 | 0.46 |
| | cis-1,2-Dichloroethylene | 70 | 50,000 | N.D. | U | 1.0 | 0.5 | 0.48 | N.D. | U | 1.0 | 0.5 | 0.48 | NA | N.D. | U | 1.0 | 0.5 | 0.48 |
| 1 | Dibromochloromethane | 0.16 | 50,000 | N.D. | U U | 1.0 | 0.5 | 0.48 | N.D. | U | 1.0 | 0.5 | 0.48 | NA | N.D. | U | 1.0 | 0.5 | 0.46 |
| 1 | Ethylbenzene | 700 | 30 | N.D. | U | 1.0 | 0.5 | 0.14 | N.D. | U | 1.0 | 0.5 | 0.14 | NA | N.D. | U | 1.0 | 0.5 | 0.14 |
| 1 | Hexachlorobutadiene | 0.86 | 6 | N.D. | U | 1.0 | 0.5 | 0.32 | N.D. | U | 1.0 | 0.5 | 0.32 | NA | N.D. | U | 1.0 | 0.5 | 0.32 |
| 1 | Methyl ethyl ketone (2-Butanone) | 7,100 | 8,400 | N.D. | U | 10 | 5.0 | 2.2 | N.D. | U | 10 | 5.0 | 2.2 | NA | N.D. | U | 10 | 5.0 | 2.2 |
| | Methyl isobutyl ketone (4-Methyl-2-Pentanone) | 2,000 | 1300 | N.D. | U | 10 | 5.0 | 4.4 | N.D. | U | 10 | 5.0 | 4.4 | NA | N.D. | U | 10 | 5.0 | 4.4 |
| 1 | Methyl tert-butyl Ether | 12 | 5 | N.D. | U | 1.0 | 0.5 | 0.31 | N.D. | U | 1.0 | 0.5 | 0.31 | NA | N.D. | U | 1.0 | 0.5 | 0.31 |
| | Methylene chloride | 4.8 | 9,100 | N.D. | U | 5.0 | 1.0 | 0.64 | N.D. | U | 5.0 | 1.0 | 0.64 | NA | N.D. | U | 5.0 | 1.0 | 0.64 |
| | Styrene | 100 | 10 | N.D. | U | 1.0 | 0.5 | 0.17 | N.D. | U | 1.0 | 0.5 | 0.17 | NA | N.D. | U | 1.0 | 0.5 | 0.17 |
| | Tetrachloroethane, 1,1,1,2- | 0.52 | 50,000 | N.D. | U | 1.0 | 0.5 | 0.40 | N.D. | U | 1.0 | 0.5 | 0.40 | NA | N.D. | U | 1.0 | 0.5 | 0.40 |
| | Tetrachloroethane, 1,1,2,2- Tetrachloroethylene | 0.067 | 500 170 | N.D. N.D. | UJ U | 1.0 | 0.5 | 0.41 | N.D. N.D. | UJ | 1.0 5.0 | 0.5 | 0.41 0.39 | NA NA | N.D. N.D. | UJ U | 1.0 5.0 | 0.5 | 0.41 0.39 |
| | Toluene | 1,000 | 40 | N.D. | U | 1.0 | 0.5 | 0.39 | N.D. | U | 1.0 | 0.5 | 0.39 | NA | N.D. | U | 1.0 | 0.5 | 0.39 |
| | trans-1,2- Dichloroethylene | 1,000 | 260 | N.D. | U | 1.0 | 0.5 | 0.24 | N.D. | U | 1.0 | 0.5 | 0.24 | NA | N.D. | U | 1.0 | 0.5 | 0.24 |
| | Trichloroethylene | 5 | 310 | N.D. | U | 1.0 | 0.5 | 0.37 | N.D. | U | 1.0 | 0.5 | 0.37 | NA | N.D. | U | 1.0 | 0.5 | 0.37 |
| | Vinyl chloride | 2 | 3,400 | N.D. | U | 1.0 | 0.5 | 0.30 | N.D. | U | 1.0 | 0.5 | 0.30 | NA | N.D. | U | 1.0 | 0.5 | 0.30 |
| | Xylenes | 10,000 | 20 | N.D. | U | 11 | 1.5 | 0.23 | N.D. | U | 11 | 1.5 | 0.23 | NA | N.D. | U | 11 | 1.5 | 0.23 |
| EPA 6020 | Dissolved Lead | 15 | 50,000 | N.D. | U | 1.0 | 0.2 | 0.0898 | N.D. | U | 1 | 0.2 | 0.0898 | NA | - | - | - | - | - |
| | n micrograms per liter (ug/L) Shaded values exceeded th | | | | | | | | | | | | | | | | | | |

The data are in micrograms per liter (µg/L). Shaded values exceeded the DOH EALs. DOH EALs DOH Tier 1 Environmental Action Levels for groundwater where groundwater is a current drinking water source and surface water is greater than 150 meters from the site (DOH, Fall 2011). DL Detection Limit or Method Detection Limit (MDL) EPA Environmental Protection Agency HD The chromatographic pattern was inconsistent with the profile of the reference fuel standard. J Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated. LOD Limit of Detection

LOQ NA N.D. Limit of Quantitation Both results for duplicate pair were non-detect, no RPD calculations Not Detected

U

Qualifiers

Q TPH-g TPH-d

Total Petroleum Hydrocarbons as gasoline Total Petroleum Hydrocarbons as diesel Undetected at DL and is reported as less than the LOD.

SECTION 4 – SUMMARY, CONCLUSIONS, AND RECOMMENDATIONS

This quarterly monitoring report presents the results of groundwater sampling conducted on July 24, 2013, at the RHSF, JBPHH, Hawaii. The RHSF is located in Halawa Heights on the Island of Oahu. The DOH Facility ID number for the RHSF is 9-102271. The DOH Release ID numbers are 990051, 010011, and 020028.

The groundwater sampling was conducted as part of the long-term groundwater and soil vapor monitoring at the RHSF, under NAVFAC Contract Number N62742-12-D-1853. The sampling was conducted in accordance with the approved WP/SAP prepared by ESI.

On July 24, 2013, ESI personnel collected groundwater samples from two monitoring wells (wells HDMW2253-03 and OWDFMW01). A summary of the analytical results is provided below.

- HDMW2253-03 Naphthalene (0.03 µg/L) was detected. The COPC was not detected at a concentration above the DOH EALs.
- OWDFMW01 TPH-d (470 and 340 μg/L), acetone (88 and 83 μg/L), naphthalene (0.081 and 0.12 μg/L), and benzene (0.42 and 0.44 μg/L) were detected in both the primary and duplicate samples. TPH-d (470 and 340 μg/L) was detected at concentrations above the DOH EALs for both drinking water toxicity and gross contamination in both samples.

Groundwater Contaminant Trends

Historical groundwater contaminant concentration trends of COPCs that exceeded the DOH EALs are presented in Appendix D. A summary of groundwater contaminant trends is provided below.

- HDMW2253-03 TPH-d was not detected during this round of quarterly sampling. TPH-d concentrations last exceeded the DOH EALs for both drinking water toxicity and gross contamination in January 2013 (600 µg/L). Naphthalene was detected but remained at a low concentration, below the DOH EALs and consistent with historical results.
- OWDFMW01 With the exception of TPH-d, groundwater contaminant concentrations remained at low concentrations and did not change significantly, or were not detected. TPH-d concentrations detected during this round of sampling were above the DOH EALs for both drinking water toxicity and gross contamination, but decreased significantly from the last event, and were the lowest they have been since July 2012.

Conclusions and Recommendations

Since the wells were last sampled (April 2013), with the exception of TPH-d, groundwater contaminant concentrations remained at low concentrations and did not change significantly,

or were not detected. TPH-d concentrations decreased in wells HDMW2253-03 and OWDFMW01. The TPH-d concentration in HDMW2253-03 was 45 μ g/L in April 2013, but was not detected during this round of sampling. TPH-d concentrations in well OWDFMW01 decreased from 1,900 μ g/L at the last round of sampling to 470 μ g/L during this round.

Based on the results of the assessment, continued groundwater monitoring at the RHSF is recommended. If the TPH-d concentrations significantly increase, the monitoring frequency should be increased to monthly, even though the two outside wells are not included in the RHSF Groundwater Protection Plan.

SECTION 5 – FUTURE WORK

GROUNDWATER SAMPLING

Future work includes the fourth quarter 2013 groundwater monitoring which is scheduled for October 2013. It is anticipated that the quarterly groundwater monitoring status report will be submitted in November 2013.

SECTION 6 – REFERENCES

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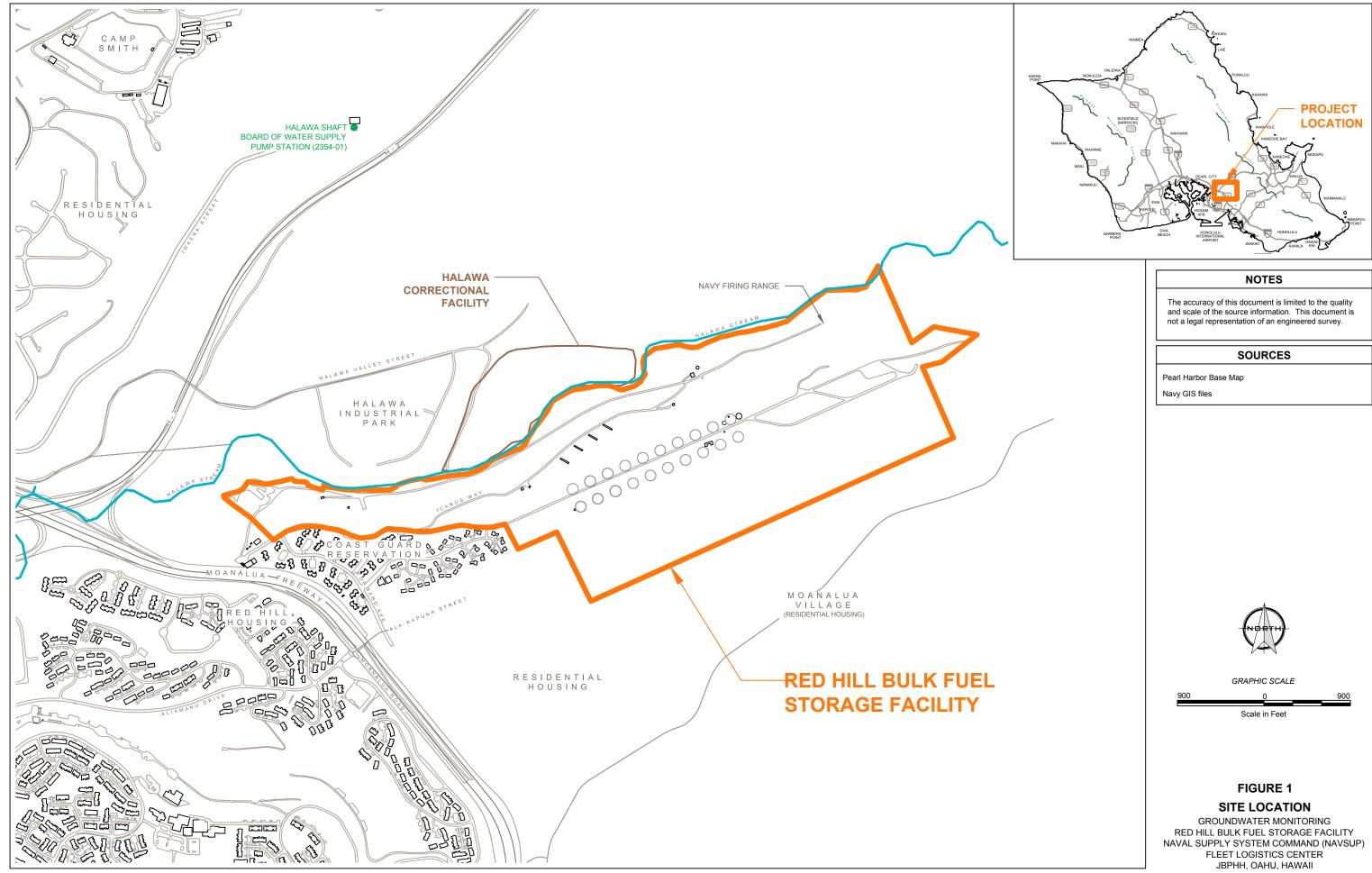
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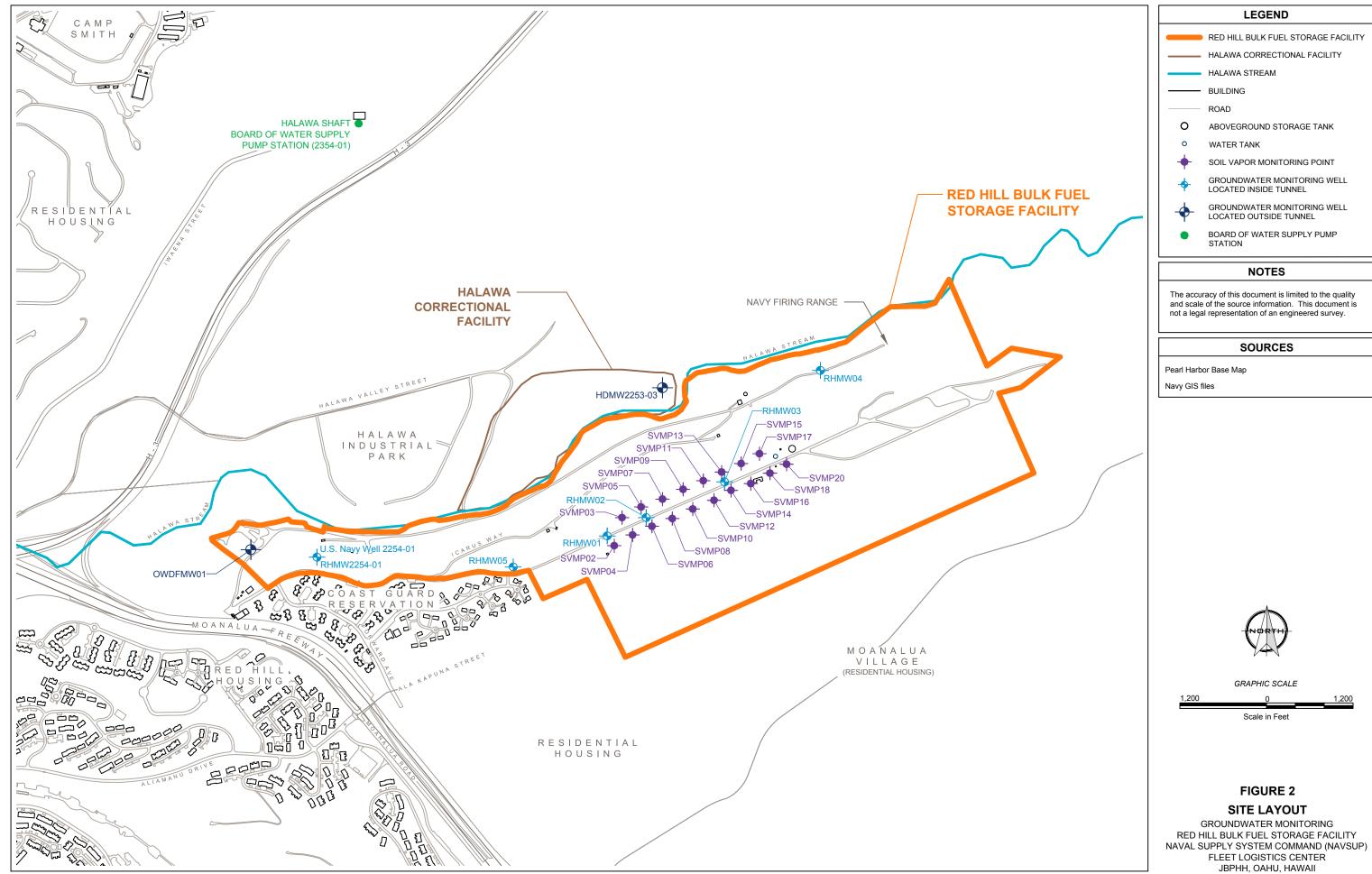
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TEC, 2010a, Quarterly Groundwater Monitoring Report – Outside (Non-Tunnel) Wells, Prepared for Navy Region Hawaii, Pearl Harbor, Hawaii, April 2010.

TEC, 2010b, Quarterly Groundwater Monitoring Report – Outside (Non-Tunnel) Wells, Prepared for Navy Region Hawaii, Pearl Harbor, Hawaii, May 2010.

FIGURES





APPENDIX A

Groundwater Sampling Logs

Groundwater Sampling Log

| Well ID: O | Well ID: OWDFMW01 Location: Red Hill Bulk Fuel St | | | I Bulk Fuel Stora | age Facility | Project No.: | 112066 |
|--------------|---|---------------------|----------------------|-------------------|--------------------|------------------|---------------------|
| Initial Wate | er Level: 12 | 20.72 ft | Date: | 7/24/2013 | | Time: 713 | 3 |
| Total Deptl | h of Well: | 145.10 ft | Personnel Involved: | | Branden | Ibara, Tina Alde | er |
| Length of S | Saturated Zon | e: | Weathe | er Conditions: | | Sunny | |
| Volume of | Water to be R | emoved: 5.0 L | Method | d of Removal: | Disposa | ble Hand Bailer | r |
| Water Leve | el After Purgin | g: <u>120.76 ft</u> | Pumpir | ng Rate: | 0.2 | 9 L/min | |
| Well Purge | e Data: | | | | | | |
| Time | Volume Removed | | nductivity mS/cm) | DO (mg/l) | Temperature | Salinity | Redox (ORP) (mV) |
| 720 | 0.0 L | 11.01 | 2.748 | 1.75 | 23.47 | <u>-</u> | -127.2 |
| 728 | 2.0 L | 10.59 | 2.724 | 1.70 | 23.60 | | -142.2 |
| 731 | 3.0 L | 10.52 | 2.751 | 1.98 | 23.60 | | -137.4 |
| 734 | 4.0 L | 10.50 | 2.765 | 1.41 | 23.60 | | -138.6 |
| 737 | 5.0 L | 10.49 | 2.765 | 1.40 | 23.60 | _ | -136.4 |
| | | · · | | | | | |
| | | · · | | | | | |
| | | | | | | | |
| | | | | | | | |
| | <u></u> | · | | | | | |
| | | · | | | | | |
| Sample W | ithdrawal Meth | nod: Dispo | sable Han | d Bailer | | | |
| Appearance | e of Sample: | | | | | | |
| | Color: | | Clear | | | | |
| | Turbidity: | | Low | | | | |
| | Sediment: | | Low | | | | |
| | Other: | | None | | | | |
| Laboratory | Analysis Para | meters and Prese | rvatives: | TPH-d - 8015; | TPH-g, VOCs - 8 | 3260; PAHs - 82 | 270c sim; |
| | - | | | lead - 6020 | | | |
| Number ar | nd Types of Sa | ample Containers: | 16 - 40ml | VOAs, 6 - 1L am | nber jar, 4 - 500m | l amber jar, 4 - | 500ml HDPE |
| Sample Ide | entification Nu | mbers: ES034 | [0815], ES | 034 MS/MSD [08 | 315]; ES035 [0900 | 0] | |
| Decontami | nation Proced | ures: Triple Rins | ed | | | | |
| Notes: YS | SI did not have | e salinity paramete | r | | | | |
| Sampled b | | Ibara, Tina Alder | | | | | |
| • | Delivered to: | Calscience Er | nvironment | | Transporters: Fee | dEx | |
| Date: 7/2 | 24/2013 | C | | | Time: <u>1100</u> | | |
| | | Cap | acity of Ca | sing (Gallons/Lir | | | |

2"-0.16• 4"-0.65 • 8"-2.61 • 10"-4.08 • 12"-5.87

Groundwater Sampling Log

| Well ID: H | DMW2253-03 | Location: | Red Hil | I Bulk Fuel Stora | age Facility | Project No.: | 112066 |
|--------------|--------------------|------------------|------------------|-------------------------------------|---------------------------------|-------------------|----------------|
| Initial Wate | er Level: | | Date: | 7/24/2013 | | Time: 913 | |
| Total Dept | h of Well: | 1575 ft | Person | nel Involved: | Branden | Ibara, Tina Aldei | |
| Length of S | Saturated Zone: | | Weathe | er Conditions: | | Sunny | |
| Volume of | Water to be Rem | noved: - | Method | d of Removal: | Disposa | able Hand Bailer | |
| Water Lev | el After Purging: | | Pumpir | ng Rate: | 0.2 | 11 L/min | |
| Well Purge | e Data: | | | | | | |
| Time | Volume | | onductivity | | Tomporatura | Solinit <i>i</i> | Redox (ORP) |
| 916 | Removed 0.0 L | рН 9.14 | (mS/cm) 0.498 | DO (mg/l) 3.33 | Temperature 23.33 | Salinity | (mV) -145.8 |
| 916 | 1.0 L | 7.19 | 0.498 | 3.02 | 23.33 | | -145.8 |
| 920 | | 6.88 | | | | | -90.2 |
| 936 | 2.0 L 3.0 L | 6.94 | 0.432 | 2.62 | 22.38 22.40 | | -90.2 |
| <u> </u> | 4.0 L | 6.77 | 0.425 | 1.96 | 22.40 | | -103.2 |
| | | | | | | | |
| Sample W | ithdrawal Method | : Disp | osable Han | d Bailer | | | |
| Appearance | ce of Sample: | | | | | | |
| | Color: | | Tan | | | | |
| | Turbidity: | | Low | | | | |
| | Sediment: | | None | | | | |
| | Other: | | None | | | | |
| Laboratory | Analysis Paramo | eters and Prese | ervatives: | <u>TPH-d - 8015;</u> lead - 6020 | TPH-g, VOCs - | 8260; PAHs - 82 | 70c sim; |
| Number ar | nd Types of Sam | ole Containers: | 6 - 40ml V | /OAs, 2 - 1L amb | oer jar, 1 - 500ml | amber jar, 1 - 25 | 0ml HDPE |
| Sample Ide | entification Numb | ers: ES036 | [0930] | | | | |
| Decontami | ination Procedure | es: Triple Rins | sed | | | | |
| Notes: YS | SI did not have sa | alinity paramete | er | | | | |
| Sampled b | | ara, Tina Alder | | | | | |
| | Delivered to: | Calscience E | nvironmenta | | Transporters: Fe | dEx | |
| Date: 7/ | 24/2013 | Car | acity of Ca | sing (Gallons/Lin | Time: <u>1100</u> hear Feet) | | |
| | | | | | | | |

APPENDIX B

Field Notes

A. Same RHAF 85 _____ Date ______3 RHSF _____ Date 7124113 Location Location Project / Client _____NAVFAC Collectell 95033 from 1045 Task: Cru Senghis RHALLOS Personnel, BI, TA Exited Tunnel 20 Propped of IDW Arrived at Fekex On-site, Health + Saloty 0710 145 Meetin 310 Gange ONDEMNOI DTW, 120.72 PID: 0.0 DTB: 145.10 Collected ESO 34, ESO 34 ns/kg ESU35 From ONDEMNOI Met DUNR at Halana Start pirge HDMW 2253-03 collected ESC36 From 6+DMW2253-03 Dropped off IDW and backed sendes Weft Reel HILL For FoelEx Left Fed Ex Back at office unleady Þ 7/24/13

APPENDIX C

Laboratory Reports



WORK ORDER NUMBER: 13-07-1752

The difference is service



AIR | SOIL | WATER | MARINE CHEMISTRY

Analytical Report For Client: Environmental Science International, Inc. Client Project Name: Red Hill LTM 112066 Attention: Robert Chong 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

Richard Villey)

Approved for release on 08/02/2013 by: Richard Villafania Project Manager

ResultLink ▶

Email your PM >



Calscience Environmental Laboratories, Inc. (Calscience) certifies that the test results provided in this report meet all NELAC requirements for parameters for which accreditation is required or available. Any exceptions to NELAC requirements are noted in the case narrative. The original report of subcontracted analyses, if any, is attached to this report. The results in this report are limited to the sample(s) tested and any reproduction thereof must be made in its entirety. The client or recipient of this report is specifically prohibited from making material changes to said report and, to the extent that such changes are made, Calscience is not responsible, legally or otherwise. The client or recipient agrees to indemnify Calscience for any defense to any litigation which may arise.



10 Lincoln Way, Garden Grove, CA 92841-1432 * TEL: (714) 895-5494 * FAX: (714) 894-7501 * www.calscience.com



Client Project Name: Red Hill LTM 112066 Work Order Number: 13-07-1752

| 1 | Work Order Narrative. | 3 |
|---|---|------------------------|
| 2 | Client Sample Data | 4 4 5 6 10 |
| 3 | Quality Control Sample Data.3.1 MS/MSD.3.2 PDS/PDSD.3.3 LCS/LCSD. | 20 20 25 26 |
| 4 | Sample Analysis Summary | 31 |
| 5 | Glossary of Terms and Qualifiers | 32 |
| 6 | Chain of Custody/Sample Receipt Form | 33 |





Work Order: 13-07-1752

Page 1 of 1

Condition Upon Receipt:

Samples were received under Chain of Custody (COC) on 07/26/13. They were assigned to Work Order 13-07-1752.

Unless otherwise noted on the Sample Receiving forms all samples were received in good condition and within the recommended EPA temperature criteria for the methods noted on the COC. The COC and Sample Receiving Documents are integral elements of the analytical report and are presented at the back of the report.

Holding Times:

All samples were analyzed within prescribed holding times (HT) and/or in accordance with the Calscience Sample Acceptance Policy unless otherwise noted in the analytical report and/or comprehensive case narrative, if required.

Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

Quality Control:

All quality control parameters (QC) were within established control limits except where noted in the QC summary forms or described further within this report.

Additional Comments:

Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are always reported on a wet weight basis.

Subcontractor Information:

Unless otherwise noted below (or on the subcontract form), no samples were subcontracted.

Return to Contents

| Environment | al Science Internationa | al, Inc. | D | ate Receiv | ved: | | | 07/26/13 |
|----------------------------------|-----------------------------|------------------------|------------------------------|----------------------|---------------|-------------------|-----------------------|-------------------|
| 354 Uluniu Street, Suite 304 | | | | Work Order: 13-07-17 | | | | |
| Kailua, HI 96 | 734-2500 | | Р | reparation | : | | | EPA 35100 |
| | | | Ν | lethod: | | | E | EPA 8015B (M |
| | | | U | nits: | | | | ug/l |
| Project: Red | Hill LTM 112066 | | | | | | Pa | age 1 of 1 |
| Client Sample N | lumber | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
| ES 034 | | 13-07-1752-2-H | 07/24/13 08:15 | Aqueous | GC 45 | 07/30/13 | 07/31/13 04:45 | 130730B09 |
| Comment(s): | - Results were evaluated to | | | the MDL (DL | _) but < RL (| LOQ), if found, a | are qualified with | a "J" flag. |
| | - TPH as DRO is quantified | d in the carbon range | e C10-C28. | | | | | |
| Parameter | | <u>Result</u> | DL | <u>LOD</u> | | LOQ | DF | <u>Qualifiers</u> |
| TPH as Diesel | | 470 | 15 | 21 | | 52 | 1.04 | HD |
| Surrogate | | Rec. (%) | Control Lir | nit <u>s</u> Qualit | fiore | | | |
| n-Octacosane | | <u>Rec. (76)</u> 90 | 51-141 | | <u>liers</u> | | | |
| n-Octacosarie | | 30 | 51-141 | | | | | |
| ES 035 | | 13-07-1752-3-Н | 07/24/13 09:00 | Aqueous | GC 45 | 07/30/13 | 07/31/13 05:02 | 130730B09 |
| Comment(s): | - Results were evaluated to | o the MDL (DL), con | centrations >= to | the MDL (DL | _) but < RL (| LOQ), if found, a | are qualified with | a "J" flag. |
| | - TPH as DRO is quantified | d in the carbon range | e C10-C28. | | | | | |
| Parameter | | <u>Result</u> | DL | LOD | | LOQ | <u>DF</u> | <u>Qualifiers</u> |
| TPH as Diesel | | 340 | 15 | 21 | | 52 | 1.04 | HD |
| Curre nete | | | Control Lin | | (| | | |
| <u>Surrogate</u> | | <u>Rec. (%)</u> 94 | <u>Control Lir</u> 51-141 | nits Qualit | liers | | | |
| n-Octacosane | | 94 | 51-141 | | | | | |
| ES 036 | | 13-07-1752-4-H | 07/24/13 09:30 | Aqueous | GC 45 | 07/30/13 | 07/31/13 05:20 | 130730B09 |
| Comment(s): | - Results were evaluated to | o the MDL (DL), con | centrations >= to | the MDL (DL |) but < RL (| LOQ), if found, a | are qualified with | a "J" flag. |
| | - TPH as DRO is quantified | d in the carbon range | e C10-C28. | | | | | |
| Parameter | | <u>Result</u> | DL | LOD | | LOQ | <u>DF</u> | <u>Qualifiers</u> |
| TPH as Diesel | | <21 | 15 | 21 | | 52 | 1.04 | U |
| Surrogate | | <u>Rec. (%)</u> | Control Lir | nits Qualit | fiers | | | |
| n-Octacosane | | 75 | 51-141 | | | | | |
| | | | | | | | | |
| Method Blank | | 099-15-516-53 | N/A | Aqueous | GC 45 | 07/30/13 | 07/31/13 01:36 | 130730B09 |
| Comment(s): | - Results were evaluated to | o the MDL (DL), con | centrations >= to | the MDL (DL | _) but < RL (| LOQ), if found, a | are qualified with | a "J" flag. |
| Parameter | | <u>Result</u> | DL | LOD | | LOQ | DF | <u>Qualifiers</u> |
| TPH as Diesel | | <20 | 15 | 20 | | 50 | 1 | U |
| | | | | | | | | |
| Curro got- | | D = - (0/) | Control 1 | | lara | | | |
| <u>Surrogate</u> n-Octacosane | | <u>Rec. (%)</u> 104 | <u>Control Lir</u> 51-141 | <u>nits</u> Qualit | fiers | | | |

Calscience nvironmental Laboratories, Inc.

| Environmental | Science Internationa | l, Inc. | | Date Receiv | ved: | | | 07/26/13 |
|----------------------------|-----------------------------|----------------------|------------------------|---------------|-----------------|------------------|-----------------------|-------------------|
| 354 Uluniu Str | eet, Suite 304 | | | Work Order: | : | | | 13-07-1752 |
| Kailua, HI 967 | 34-2500 | | | Preparation | | | EF | PA 3020A Total |
| - | | | | Method: | | | | EPA 6020 |
| | | | | Units: | | | | ug/L |
| Project: Red H | iill LTM 112066 | | | | | | Р | age 1 of 1 |
| Client Sample Nu | mber | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
| ES 034 | | 13-07-1752-2-G | 07/24/13 08:15 | Aqueous | ICP/MS 04 | 07/31/13 | 07/31/13 17:51 | 130731L02D |
| Comment(s): | - Results were evaluated to | the MDL (DL), con | centrations >= t | o the MDL (DL |) but < RL (LOC | Q), if found, a | are qualified with | a "J" flag. |
| Parameter | | <u>Result</u> | DL | LOD | LOC | <u>2</u> | DF | <u>Qualifiers</u> |
| Lead | | <0.200 | 0.0898 | 0.200 | 1.00 |) | 1 | U |
| ES 035 | | 13-07-1752-3-G | 07/24/13 09:00 | Aqueous | ICP/MS 04 | 07/31/13 | 07/31/13 17:55 | 130731L02D |
| Comment(s): | - Results were evaluated to | the MDL (DL), con | centrations >= t | o the MDL (DL |) but < RL (LOC | Q), if found, a | are qualified with | a "J" flag. |
| Parameter erementer | | <u>Result</u> | <u>DL</u> | LOD | LOC | 2 | DF | <u>Qualifiers</u> |
| Lead | | <0.200 | 0.0898 | 0.200 | 1.00 |) | 1 | U |
| ES 036 | | 13-07-1752-4-G | 07/24/13 09:30 | Aqueous | ICP/MS 04 | 07/31/13 | 07/31/13 17:59 | 130731L02D |
| Comment(s): | - Results were evaluated to | the MDL (DL), con | centrations >= t | o the MDL (DL |) but < RL (LOC | Q), if found, a | are qualified with | a "J" flag. |
| Parameter | | <u>Result</u> | DL | LOD | <u>LOC</u> | <u>2</u> | <u>DF</u> | <u>Qualifiers</u> |
| Lead | | <0.200 | 0.0898 | 0.200 | 1.00 |) | 1 | U |
| Method Blank | | 099-14-497-42 | N/A | Aqueous | ICP/MS 04 | 07/31/13 | 07/31/13 17:23 | 130731L02D |
| Comment(s): | - Results were evaluated to | the MDL (DL), con | centrations >= t | o the MDL (DL |) but < RL (LOC | Q), if found, a | are qualified with | a "J" flag. |
| D / | | Result | DL | LOD | LOC | ` | DF | Qualifiers |
| Parameter | | Result | | LOD | <u>LU(</u> | <u>x</u> | | Quaimers |

Analytical Report



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

| Date Received: | 07/26/13 |
|----------------|--------------------|
| Work Order: | 13-07-1752 |
| Preparation: | EPA 3510C |
| Method: | EPA 8270C SIM PAHs |
| Units: | ug/L |
| | Page 1 of 4 |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--------------------------------------|----------------------|------------------------|-------------------|-----------------|------------------|-----------------------|-------------------|
| ES 034 | 13-07-1752-2-l | 07/24/13 08:15 | Aqueous | GC/MS AAA | 07/29/13 | 08/01/13 13:55 | 130729L12 |
| Comment(s): - Results were evaluated | to the MDL (DL), con | centrations >= to th | he MDL (DL |) but < RL (LOC |), if found, a | are qualified with | a "J" flag. |
| Parameter | <u>Result</u> | DL | LOD | LOC | 2 | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | 0.081 | 0.023 | 0.050 | 0.20 |) | 0.995 | J |
| 2-Methylnaphthalene | <0.050 | 0.026 | 0.050 | 0.20 |) | 0.995 | U |
| 1-Methylnaphthalene | <0.050 | 0.028 | 0.050 | 0.20 |) | 0.995 | U |
| Acenaphthylene | <0.050 | 0.018 | 0.050 | 0.20 |) | 0.995 | U |
| Acenaphthene | <0.050 | 0.021 | 0.050 | 0.20 |) | 0.995 | U |
| Fluorene | <0.050 | 0.024 | 0.050 | 0.20 |) | 0.995 | U |
| Phenanthrene | <0.050 | 0.030 | 0.050 | 0.20 |) | 0.995 | U |
| Anthracene | <0.050 | 0.034 | 0.050 | 0.20 |) | 0.995 | U |
| Fluoranthene | <0.050 | 0.027 | 0.050 | 0.20 |) | 0.995 | U |
| Pyrene | <0.050 | 0.025 | 0.050 | 0.20 |) | 0.995 | U |
| Benzo (a) Anthracene | <0.050 | 0.024 | 0.050 | 0.20 |) | 0.995 | U |
| Chrysene | <0.050 | 0.019 | 0.050 | 0.20 |) | 0.995 | U |
| Benzo (k) Fluoranthene | <0.050 | 0.023 | 0.050 | 0.20 |) | 0.995 | U |
| Benzo (b) Fluoranthene | <0.050 | 0.025 | 0.050 | 0.20 |) | 0.995 | U |
| Benzo (a) Pyrene | <0.050 | 0.036 | 0.050 | 0.20 |) | 0.995 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.050 | 0.022 | 0.050 | 0.20 |) | 0.995 | U |
| Dibenz (a,h) Anthracene | <0.050 | 0.027 | 0.050 | 0.20 |) | 0.995 | U |
| Benzo (g,h,i) Perylene | <0.050 | 0.022 | 0.050 | 0.20 |) | 0.995 | U |
| Surrogate | <u>Rec. (%)</u> | Control Limi | <u>ts Qualifi</u> | iers | | | |
| Nitrobenzene-d5 | 64 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 63 | 33-144 | | | | | |
| p-Terphenyl-d14 | 76 | 23-160 | | | | | |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

| Date Received: | 07/26/13 |
|----------------|--------------------|
| Work Order: | 13-07-1752 |
| Preparation: | EPA 3510C |
| Method: | EPA 8270C SIM PAHs |
| Units: | ug/L |
| | Page 2 of 4 |

Project: Red Hill LTM 112066

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--|----------------------|------------------------|-------------------|------------------|------------------|-----------------------|-------------------|
| ES 035 | 13-07-1752-3-l | 07/24/13 09:00 | Aqueous | GC/MS AAA | 07/29/13 | 08/01/13 14:21 | 130729L12 |
| Comment(s): - Results were evaluated t | o the MDL (DL), con | centrations >= to t | he MDL (DL | .) but < RL (LOC | Q), if found, a | are qualified with a | a "J" flag. |
| Parameter | <u>Result</u> | <u>DL</u> | LOD | LOC | <u>2</u> | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | 0.12 | 0.022 | 0.048 | 0.19 |) | 0.965 | J |
| 2-Methylnaphthalene | <0.048 | 0.026 | 0.048 | 0.19 |) | 0.965 | U |
| 1-Methylnaphthalene | <0.048 | 0.027 | 0.048 | 0.19 |) | 0.965 | U |
| Acenaphthylene | <0.048 | 0.017 | 0.048 | 0.19 |) | 0.965 | U |
| Acenaphthene | <0.048 | 0.020 | 0.048 | 0.19 |) | 0.965 | U |
| Fluorene | <0.048 | 0.024 | 0.048 | 0.19 |) | 0.965 | U |
| Phenanthrene | <0.048 | 0.029 | 0.048 | 0.19 |) | 0.965 | U |
| Anthracene | <0.048 | 0.033 | 0.048 | 0.19 |) | 0.965 | U |
| Fluoranthene | <0.048 | 0.026 | 0.048 | 0.19 |) | 0.965 | U |
| Pyrene | <0.048 | 0.024 | 0.048 | 0.19 |) | 0.965 | U |
| Benzo (a) Anthracene | <0.048 | 0.023 | 0.048 | 0.19 |) | 0.965 | U |
| Chrysene | <0.048 | 0.018 | 0.048 | 0.19 |) | 0.965 | U |
| Benzo (k) Fluoranthene | <0.048 | 0.023 | 0.048 | 0.19 |) | 0.965 | U |
| Benzo (b) Fluoranthene | <0.048 | 0.024 | 0.048 | 0.19 |) | 0.965 | U |
| Benzo (a) Pyrene | <0.048 | 0.035 | 0.048 | 0.19 |) | 0.965 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.048 | 0.021 | 0.048 | 0.19 |) | 0.965 | U |
| Dibenz (a,h) Anthracene | <0.048 | 0.026 | 0.048 | 0.19 |) | 0.965 | U |
| Benzo (g,h,i) Perylene | <0.048 | 0.021 | 0.048 | 0.19 |) | 0.965 | U |
| Surrogate | <u>Rec. (%)</u> | Control Lim | <u>its Qualif</u> | iers | | | |
| Nitrobenzene-d5 | 75 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 70 | 33-144 | | | | | |
| p-Terphenyl-d14 | 80 | 23-160 | | | | | |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

| Date Received: | 07/26/13 |
|----------------|--------------------|
| Work Order: | 13-07-1752 |
| Preparation: | EPA 3510C |
| Method: | EPA 8270C SIM PAHs |
| Units: | ug/L |
| | Page 3 of 4 |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--|----------------------|------------------------|-------------------|------------------|------------------|-----------------------|-------------|
| ES 036 | 13-07-1752-4-I | 07/24/13 09:30 | Aqueous | GC/MS AAA | 07/29/13 | 08/01/13 14:47 | 130729L12 |
| Comment(s): - Results were evaluated t | o the MDL (DL), con | centrations >= to t | he MDL (DL | .) but < RL (LOC | Q), if found, a | are qualified with a | a "J" flag. |
| Parameter | <u>Result</u> | DL | LOD | LOC | 2 | <u>DF</u> | Qualifiers |
| Naphthalene | 0.030 | 0.022 | 0.047 | 0.19 |) | 0.945 | J |
| 2-Methylnaphthalene | <0.047 | 0.025 | 0.047 | 0.19 |) | 0.945 | U |
| 1-Methylnaphthalene | <0.047 | 0.027 | 0.047 | 0.19 |) | 0.945 | U |
| Acenaphthylene | <0.047 | 0.017 | 0.047 | 0.19 |) | 0.945 | U |
| Acenaphthene | <0.047 | 0.020 | 0.047 | 0.19 |) | 0.945 | U |
| Fluorene | <0.047 | 0.023 | 0.047 | 0.19 |) | 0.945 | U |
| Phenanthrene | <0.047 | 0.029 | 0.047 | 0.19 |) | 0.945 | U |
| Anthracene | <0.047 | 0.032 | 0.047 | 0.19 |) | 0.945 | U |
| Fluoranthene | <0.047 | 0.026 | 0.047 | 0.19 |) | 0.945 | U |
| Pyrene | <0.047 | 0.023 | 0.047 | 0.19 |) | 0.945 | U |
| Benzo (a) Anthracene | <0.047 | 0.022 | 0.047 | 0.19 |) | 0.945 | U |
| Chrysene | <0.047 | 0.018 | 0.047 | 0.19 |) | 0.945 | U |
| Benzo (k) Fluoranthene | <0.047 | 0.022 | 0.047 | 0.19 |) | 0.945 | U |
| Benzo (b) Fluoranthene | <0.047 | 0.024 | 0.047 | 0.19 |) | 0.945 | U |
| Benzo (a) Pyrene | <0.047 | 0.034 | 0.047 | 0.19 |) | 0.945 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.047 | 0.021 | 0.047 | 0.19 |) | 0.945 | U |
| Dibenz (a,h) Anthracene | <0.047 | 0.025 | 0.047 | 0.19 |) | 0.945 | U |
| Benzo (g,h,i) Perylene | <0.047 | 0.021 | 0.047 | 0.19 |) | 0.945 | U |
| Surrogate | <u>Rec. (%)</u> | Control Limi | <u>its</u> Qualif | iers | | | |
| Nitrobenzene-d5 | 68 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 66 | 33-144 | | | | | |
| p-Terphenyl-d14 | 72 | 23-160 | | | | | |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

| Date Received: | 07/26/13 |
|----------------|--------------------|
| Work Order: | 13-07-1752 |
| Preparation: | EPA 3510C |
| Method: | EPA 8270C SIM PAHs |
| Units: | ug/L |
| | Page 4 of 4 |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|---|----------------------|------------------------|--------------------------|------------------|------------------|-----------------------|-------------------|
| Method Blank | 099-15-148-17 | N/A | Aqueous | GC/MS AAA | 07/29/13 | 08/01/13 13:28 | 130729L12 |
| Comment(s): - Results were evaluated to | the MDL (DL), con | centrations >= to t | he MDL (DL | .) but < RL (LOC | Q), if found, a | re qualified with a | a "J" flag. |
| Parameter | <u>Result</u> | <u>DL</u> | <u>LOD</u> | LOC | 2 | <u>DF</u> | <u>Qualifiers</u> |
| Naphthalene | <0.051 | 0.023 | 0.051 | 0.20 |) | 1.02 | U |
| 2-Methylnaphthalene | <0.051 | 0.027 | 0.051 | 0.20 |) | 1.02 | U |
| 1-Methylnaphthalene | <0.051 | 0.029 | 0.051 | 0.20 |) | 1.02 | U |
| Acenaphthylene | <0.051 | 0.018 | 0.051 | 0.20 |) | 1.02 | U |
| Acenaphthene | <0.051 | 0.021 | 0.051 | 0.20 |) | 1.02 | U |
| Fluorene | <0.051 | 0.025 | 0.051 | 0.20 |) | 1.02 | U |
| Phenanthrene | <0.051 | 0.031 | 0.051 | 0.20 |) | 1.02 | U |
| Anthracene | <0.051 | 0.035 | 0.051 | 0.20 |) | 1.02 | U |
| Fluoranthene | <0.051 | 0.028 | 0.051 | 0.20 |) | 1.02 | U |
| Pyrene | <0.051 | 0.025 | 0.051 | 0.20 |) | 1.02 | U |
| Benzo (a) Anthracene | <0.051 | 0.024 | 0.051 | 0.20 |) | 1.02 | U |
| Chrysene | <0.051 | 0.019 | 0.051 | 0.20 |) | 1.02 | U |
| Benzo (k) Fluoranthene | <0.051 | 0.024 | 0.051 | 0.20 |) | 1.02 | U |
| Benzo (b) Fluoranthene | <0.051 | 0.025 | 0.051 | 0.20 |) | 1.02 | U |
| Benzo (a) Pyrene | <0.051 | 0.037 | 0.051 | 0.20 |) | 1.02 | U |
| Indeno (1,2,3-c,d) Pyrene | <0.051 | 0.022 | 0.051 | 0.20 |) | 1.02 | U |
| Dibenz (a,h) Anthracene | <0.051 | 0.027 | 0.051 | 0.20 |) | 1.02 | U |
| Benzo (g,h,i) Perylene | <0.051 | 0.022 | 0.051 | 0.20 |) | 1.02 | U |
| Surrogate | <u>Rec. (%)</u> | Control Limi | <u>its</u> <u>Qualif</u> | iers | | | |
| Nitrobenzene-d5 | 74 | 28-139 | | | | | |
| 2-Fluorobiphenyl | 70 | 33-144 | | | | | |
| p-Terphenyl-d14 | 80 | 23-160 | | | | | |



Units:

Environmental Science International, Inc.Date Received:354 Uluniu Street, Suite 304Work Order:Kailua, HI 96734-2500Preparation:Method:Method:

| 07/26/13 |
|-------------------|
| 13-07-1752 |
| EPA 5030C |
| GC/MS / EPA 8260B |
| ug/L |
| Page 1 of 10 |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID | | | |
|--|----------------------|------------------------|---------|------------|------------------|-----------------------|-------------------|--|--|--|
| ES Trip | 13-07-1752-1-A | 07/24/13 07:00 | Aqueous | GC/MS LL | 07/30/13 | 07/31/13 01:23 | 130730L02 | | | |
| Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. | | | | | | | | | | |
| Parameter | <u>Result</u> | DL | LOD | LOC | <u>ם ב</u> | <u>F</u> | Qualifiers | | | |
| Acetone | <10 | 6.0 | 10 | 20 | 1 | | U | | | |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | | U | | | |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1 | | U | | | |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1 | | U | | | |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1 | | U | | | |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1 | | U | | | |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1 | | U | | | |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1 | | U | | | |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1 | | U | | | |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1 | | U | | | |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1 | | U | | | |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1 | | U | | | |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1 | | U | | | |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U | | | |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1 | | U | | | |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U | | | |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1 | | U | | | |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1 | | U | | | |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | | U | | | |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1 | | U | | | |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1 | | U | | | |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1 | | U | | | |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | | | U | | | |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | | | U | | | |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1 | | U | | | |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1 | | U | | | |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | | | U | | | |



| Environmental Science International, Inc | Date | Received: | | 07/26/13 | | |
|--|-----------------|----------------|------------|------------|----|-------------------|
| 354 Uluniu Street, Suite 304 | Work | Order: | | 13-07-1752 | | |
| Kailua, HI 96734-2500 | | Prepa | aration: | | | EPA 5030C |
| | | Meth | od: | | | GC/MS / EPA 8260B |
| | | Units | : | | | ug/L |
| Project: Red Hill LTM 112066 | | | | | | Page 2 of 10 |
| Parameter | <u>Result</u> | DL | LOD | LOQ | DF | Qualifiers |
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1 | U |
| Gasoline Range Organics | <30 | 13 | 30 | 50 | 1 | U |
| Surrogate | <u>Rec. (%)</u> | Control Limits | Qualifiers | | | |
| Dibromofluoromethane | 94 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 90 | 80-134 | | | | |
| Toluene-d8 | 99 | 80-120 | | | | |
| Toluene-d8-TPPH | 95 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 92 | 80-120 | | | | |

Page 3 of 10



| Environmental Science International, Inc. | Date Received: | 07/26/13 |
|---|----------------|-------------------|
| 354 Uluniu Street, Suite 304 | Work Order: | 13-07-1752 |
| Kailua, HI 96734-2500 | Preparation: | EPA 5030C |
| | Method: | GC/MS / EPA 8260B |
| | Units: | ug/L |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID | | | | |
|--|----------------------|------------------------|---------|------------|------------------|-----------------------|-------------------|--|--|--|--|
| ES 034 | 13-07-1752-2-A | 07/24/13 08:15 | Aqueous | GC/MS LL | 07/30/13 | 07/31/13 01:51 | 130730L02 | | | | |
| Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. | | | | | | | | | | | |
| Parameter | <u>Result</u> | DL | LOD | LOC | <u>ם ב</u> | <u>IF</u> | <u>Qualifiers</u> | | | | |
| Acetone | 88 | 6.0 | 10 | 20 | 1 | | | | | | |
| Benzene | 0.42 | 0.14 | 0.50 | 1.0 | 1 | | J | | | | |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1 | | U | | | | |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1 | | U | | | | |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1 | | U | | | | |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1 | | U | | | | |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1 | | U | | | | |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1 | | U | | | | |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1 | | U | | | | |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1 | | U | | | | |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1 | | U | | | | |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1 | | U | | | | |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1 | | U | | | | |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U | | | | |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1 | | U | | | | |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1 | | U | | | | |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | | |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | | |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | | U | | | | |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1 | | U | | | | |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1 | | U | | | | |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1 | | U | | | | |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1 | | U | | | | |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1 | | U | | | | |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1 | | U | | | | |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1 | | U | | | | |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1 | | U | | | | |



| Environmental Science International, Inc |). | Date | Received: | | 07/26/13 | | |
|--|-----------------|----------------|-------------------|-----|------------|-------------------|--|
| 354 Uluniu Street, Suite 304 | Work | Order: | | | 13-07-1752 | | |
| Kailua, HI 96734-2500 | | Prepa | aration: | | | EPA 5030C | |
| | | Meth | od: | | | GC/MS / EPA 8260B | |
| | | Units | : | | | ug/L | |
| Project: Red Hill LTM 112066 | | | | | | Page 4 of 10 | |
| Parameter | <u>Result</u> | DL | LOD | LOQ | DF | Qualifiers | |
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1 | U | |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | U | |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1 | U | |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1 | U | |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1 | U | |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1 | U | |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1 | U | |
| Gasoline Range Organics | <30 | 13 | 30 | 50 | 1 | U | |
| Surrogate | <u>Rec. (%)</u> | Control Limits | <u>Qualifiers</u> | | | | |
| Dibromofluoromethane | 93 | 80-126 | | | | | |
| 1,2-Dichloroethane-d4 | 90 | 80-134 | | | | | |
| Toluene-d8 | 98 | 80-120 | | | | | |
| Toluene-d8-TPPH | 94 | 88-112 | | | | | |
| 1,4-Bromofluorobenzene | 92 | 80-120 | | | | | |



Environmental Science International, Inc.Date Received:354 Uluniu Street, Suite 304Work Order:Kailua, HI 96734-2500Preparation:Method:Units:

EPA 5030C GC/MS / EPA 8260B ug/L

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07/26/13

13-07-1752

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID | | | |
|--|----------------------|------------------------|---------|------------|--------------------|-----------------------|-------------------|--|--|--|
| ES 035 | 13-07-1752-3-A | 07/24/13 09:00 | Aqueous | GC/MS LL | 07/30/13 | 07/31/13 02:18 | 130730L02 | | | |
| Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. | | | | | | | | | | |
| Parameter | <u>Result</u> | DL | LOD | LOC | <u>2</u> <u>DI</u> | <u>F</u> | <u>Qualifiers</u> | | | |
| Acetone | 83 | 6.0 | 10 | 20 | 1 | | | | | |
| Benzene | 0.44 | 0.14 | 0.50 | 1.0 | 1 | | J | | | |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1 | | U | | | |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1 | | U | | | |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1 | | U | | | |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1 | | U | | | |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1 | | U | | | |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1 | | U | | | |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1 | | U | | | |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1 | | U | | | |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1 | | U | | | |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1 | | U | | | |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1 | | U | | | |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U | | | |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1 | | U | | | |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U | | | |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1 | | U | | | |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1 | | U | | | |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | | U | | | |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1 | | U | | | |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1 | | U | | | |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1 | | U | | | |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1 | | U | | | |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1 | | U | | | |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1 | | U | | | |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1 | | U | | | |



| Environmental Science International, Inc. | | | Received: | | 07/26/13 | |
|---|-----------------|----------------|-------------------|------------|----------|-------------------|
| 354 Uluniu Street, Suite 304 | Work | Order: | | 13-07-1752 | | |
| Kailua, HI 96734-2500 | | Prepa | aration: | | | EPA 5030C |
| | | Meth | od: | | | GC/MS / EPA 8260B |
| | | Units | : | | | ug/L |
| Project: Red Hill LTM 112066 | | | | | | Page 6 of 10 |
| Parameter | <u>Result</u> | <u>DL</u> | LOD | LOQ | DF | Qualifiers |
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1 | U |
| Gasoline Range Organics | <30 | 13 | 30 | 50 | 1 | U |
| Surrogate | <u>Rec. (%)</u> | Control Limits | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 93 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 93 | 80-134 | | | | |
| Toluene-d8 | 101 | 80-120 | | | | |
| Toluene-d8-TPPH | 97 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 92 | 80-120 | | | | |



Environmental Science International, Inc. 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500

| Date Received: | 07/26/13 |
|----------------|-------------------|
| Work Order: | 13-07-1752 |
| Preparation: | EPA 5030C |
| Method: | GC/MS / EPA 8260B |
| Units: | ug/L |
| | Page 7 of 10 |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID | | | |
|--|----------------------|------------------------|---------|------------|-------------------|-----------------------|-------------------|--|--|--|
| ES 036 | 13-07-1752-4-A | 07/24/13 09:30 | Aqueous | GC/MS LL | 07/30/13 | 07/31/13 02:46 | 130730L02 | | | |
| Comment(s): - Results were evaluated to the MDL (DL), concentrations >= to the MDL (DL) but < RL (LOQ), if found, are qualified with a "J" flag. | | | | | | | | | | |
| Parameter | <u>Result</u> | DL | LOD | LOC | <u>2</u> <u>D</u> | <u>F</u> | <u>Qualifiers</u> | | | |
| Acetone | <10 | 6.0 | 10 | 20 | 1 | | U | | | |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | | U | | | |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1 | | U | | | |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1 | | U | | | |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1 | | U | | | |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1 | | U | | | |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1 | | U | | | |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1 | | U | | | |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1 | | U | | | |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1 | | U | | | |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1 | | U | | | |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1 | | U | | | |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1 | | U | | | |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U | | | |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1 | | U | | | |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U | | | |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1 | | U | | | |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1 | | U | | | |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U | | | |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | | U | | | |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1 | | U | | | |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1 | | U | | | |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U | | | |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1 | | U | | | |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1 | | U | | | |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1 | | U | | | |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1 | | U | | | |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1 | | U | | | |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1 | | U | | | |



| Environmental Science International 354 Uluniu Street, Suite 304 Kailua, HI 96734-2500 | Work | Received: Order: aration: od: | | 07/26/13 13-07-1752 EPA 5030C GC/MS / EPA 8260B | | |
|--|-----------------|--|-------------------|--|----|----------------------|
| Project: Red Hill LTM 112066 | | Units | : | | | ug/L Page 8 of 10 |
| | | | | | | |
| Parameter | <u>Result</u> | DL | LOD | LOQ | DF | Qualifiers |
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1 | U |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | U |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1 | U |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1 | U |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1 | U |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1 | U |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1 | U |
| Gasoline Range Organics | <30 | 13 | 30 | 50 | 1 | U |
| Surrogate | <u>Rec. (%)</u> | Control Limits | <u>Qualifiers</u> | | | |
| Dibromofluoromethane | 97 | 80-126 | | | | |
| 1,2-Dichloroethane-d4 | 95 | 80-134 | | | | |
| Toluene-d8 | 99 | 80-120 | | | | |
| Toluene-d8-TPPH | 96 | 88-112 | | | | |
| 1,4-Bromofluorobenzene | 89 | 80-120 | | | | |
| | | | | | | |

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| Environmental Science International, Inc. | |
|---|--|
| 354 Uluniu Street, Suite 304 | |
| Kailua, HI 96734-2500 | |

| Date Received: | 07/26/13 |
|----------------|-------------------|
| Work Order: | 13-07-1752 |
| Preparation: | EPA 5030C |
| Method: | GC/MS / EPA 8260B |
| Units: | ug/L |
| | Page 9 of 10 |

| Client Sample Number | Lab Sample Number | Date/Time Collected | Matrix | Instrument | Date Prepared | Date/Time Analyzed | QC Batch ID |
|--|----------------------|------------------------|---------------|------------------|-------------------|-----------------------|-------------------|
| Method Blank | 099-13-057-26 | N/A | Aqueous | GC/MS LL | 07/30/13 | 07/30/13 20:20 | 130730L02 |
| Comment(s): - Results were evaluated t | the MDL (DL), cor | centrations >= t | o the MDL (DL | _) but < RL (LOC | Q), if found, are | qualified with | a "J" flag. |
| Parameter | <u>Result</u> | DL | LOD | LOC | <u>2</u> <u>D</u> | <u>)F</u> | <u>Qualifiers</u> |
| Acetone | <10 | 6.0 | 10 | 20 | 1 | | U |
| Benzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | | U |
| Bromodichloromethane | <0.50 | 0.21 | 0.50 | 5.0 | 1 | | U |
| Bromoform | <1.0 | 0.50 | 1.0 | 10 | 1 | | U |
| Bromomethane | <5.0 | 3.9 | 5.0 | 20 | 1 | | U |
| 2-Butanone | <5.0 | 2.2 | 5.0 | 10 | 1 | | U |
| Carbon Tetrachloride | <0.50 | 0.23 | 0.50 | 1.0 | 1 | | U |
| Chlorobenzene | <0.50 | 0.17 | 0.50 | 5.0 | 1 | | U |
| Chloroethane | <5.0 | 2.3 | 5.0 | 10 | 1 | | U |
| Chloroform | <0.50 | 0.46 | 0.50 | 5.0 | 1 | | U |
| Chloromethane | <2.0 | 1.8 | 2.0 | 10 | 1 | | U |
| Dibromochloromethane | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U |
| 1,2-Dibromo-3-Chloropropane | <2.0 | 1.2 | 2.0 | 10 | 1 | | U |
| 1,2-Dibromoethane | <0.50 | 0.36 | 0.50 | 1.0 | 1 | | U |
| 1,2-Dichlorobenzene | <0.50 | 0.46 | 0.50 | 1.0 | 1 | | U |
| 1,3-Dichlorobenzene | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U |
| 1,4-Dichlorobenzene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U |
| 1,1-Dichloroethane | <0.50 | 0.28 | 0.50 | 5.0 | 1 | | U |
| 1,2-Dichloroethane | <0.50 | 0.24 | 0.50 | 1.0 | 1 | | U |
| 1,1-Dichloroethene | <0.50 | 0.43 | 0.50 | 1.0 | 1 | | U |
| c-1,2-Dichloroethene | <0.50 | 0.48 | 0.50 | 1.0 | 1 | | U |
| t-1,2-Dichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | | U |
| 1,2-Dichloropropane | <0.50 | 0.42 | 0.50 | 5.0 | 1 | | U |
| c-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U |
| t-1,3-Dichloropropene | <0.50 | 0.25 | 0.50 | 1.0 | 1 | | U |
| Ethylbenzene | <0.50 | 0.14 | 0.50 | 1.0 | 1 | | U |
| Methylene Chloride | <1.0 | 0.64 | 1.0 | 5.0 | 1 | | U |
| 4-Methyl-2-Pentanone | <5.0 | 4.4 | 5.0 | 10 | 1 | | U |
| Styrene | <0.50 | 0.17 | 0.50 | 1.0 | 1 | | U |
| 1,1,1,2-Tetrachloroethane | <0.50 | 0.40 | 0.50 | 1.0 | 1 | | U |
| 1,1,2,2-Tetrachloroethane | <0.50 | 0.41 | 0.50 | 1.0 | 1 | | U |
| Tetrachloroethene | <0.50 | 0.39 | 0.50 | 5.0 | 1 | | U |
| Toluene | <0.50 | 0.24 | 0.50 | 1.0 | 1 | | U |
| 1,2,4-Trichlorobenzene | <1.0 | 0.50 | 1.0 | 5.0 | 1 | | U |
| 1,1,1-Trichloroethane | <0.50 | 0.30 | 0.50 | 5.0 | 1 | | U |
| Hexachloro-1,3-Butadiene | <0.50 | 0.32 | 0.50 | 1.0 | 1 | | U |

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| Environmental Science International, In | С. | Date | Received: | | 07/26/13 | | | |
|---|-----------------|----------------|-------------------|-----|-----------|-------------------|--|--|
| 354 Uluniu Street, Suite 304 | | Work | Order: | | | 13-07-1752 | | |
| Kailua, HI 96734-2500 | Prepa | aration: | | | EPA 5030C | | | |
| Method: | | | | | | GC/MS / EPA 8260B | | |
| | | Units | : | | | ug/L | | |
| Project: Red Hill LTM 112066 | | | | | | Page 10 of 10 | | |
| Parameter | <u>Result</u> | DL | LOD | LOQ | DF | Qualifiers | | |
| 1,1,2-Trichloroethane | <0.50 | 0.38 | 0.50 | 1.0 | 1 | U | | |
| Trichloroethene | <0.50 | 0.37 | 0.50 | 1.0 | 1 | U | | |
| 1,2,3-Trichloropropane | <1.0 | 0.64 | 1.0 | 5.0 | 1 | U | | |
| Vinyl Chloride | <0.50 | 0.30 | 0.50 | 1.0 | 1 | U | | |
| p/m-Xylene | <1.0 | 0.30 | 1.0 | 10 | 1 | U | | |
| o-Xylene | <0.50 | 0.23 | 0.50 | 1.0 | 1 | U | | |
| Methyl-t-Butyl Ether (MTBE) | <0.50 | 0.31 | 0.50 | 1.0 | 1 | U | | |
| Gasoline Range Organics | <30 | 13 | 30 | 50 | 1 | U | | |
| Surrogate | <u>Rec. (%)</u> | Control Limits | <u>Qualifiers</u> | | | | | |
| Dibromofluoromethane | 91 | 80-126 | | | | | | |
| 1,2-Dichloroethane-d4 | 88 | 80-134 | | | | | | |
| Toluene-d8 | 99 | 80-120 | | | | | | |
| Toluene-d8-TPPH | 95 | 88-112 | | | | | | |
| 1,4-Bromofluorobenzene | 92 | 80-120 | | | | | | |



Quality Control - Spike/Spike Duplicate

| Environmental Science International, Inc. | Date Received: | 07/26/13 |
|---|----------------|---------------|
| 354 Uluniu Street, Suite 304 | Work Order: | 13-07-1752 |
| Kailua, HI 96734-2500 | Preparation: | EPA 3510C |
| | Method: | EPA 8015B (M) |
| Project: Red Hill LTM 112066 | | Page 1 of 5 |

Project: Red Hill LTM 112066

| Quality Control Sample ID | | | Matrix | | Date Prepared | | Date Analyzed | MS/MSD Batch Numbe | | Number |
|---------------------------|-------------------------------|-----------------------|--------------------|--------------------|---------------------|----------------------------|-----------------|--------------------|---------------|-------------------|
| ES 034 | | Aqueou | ıs | GC 45 | 07/30/ | 13 | 07/31/13 03:01 | 130 | 730S09 | |
| Parameter | <u>Sample</u> <u>Conc.</u> | <u>Spike</u> Added | <u>MS</u> Conc. | <u>MS</u> %Rec. | <u>MSD</u> Conc. | <u>MSD</u> <u>%Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| TPH as Diesel | 471.1 | 4000 | 4828 | 109 | 5043 | 114 | 55-133 | 4 | 0-30 | |

RPD: Relative Percent Difference. CL: Control Limits



Quality Control - Spike/Spike Duplicate

| Environmental Science International, Inc. | Date Received: | 07/26/13 |
|---|----------------|-----------------|
| 354 Uluniu Street, Suite 304 | Work Order: | 13-07-1752 |
| Kailua, HI 96734-2500 | Preparation: | EPA 3020A Total |
| | Method: | EPA 6020 |
| Project: Red Hill LTM 112066 | | Page 2 of 5 |

| Quality Control Sample ID | | Matrix | | Instrument | Date P | repared | Date Analyzed | MS | /MSD Batch | n Number |
|---------------------------|-------------------------------|-----------------------|--------------------|--------------------|--------------|---------------------|-----------------|------------|------------|-------------------|
| ES 034 | | Aqueo | us | ICP/MS 04 | 07/31/ | 13 | 07/31/13 17:35 | 130 | 731S02 | |
| Parameter | <u>Sample</u> <u>Conc.</u> | <u>Spike</u> Added | <u>MS</u> Conc. | <u>MS</u> %Rec. | MSD Conc. | <u>MSD</u> %Rec. | <u>%Rec. CL</u> | <u>RPD</u> | RPD CL | <u>Qualifiers</u> |
| l ead | ND | 100.0 | 109.5 | 109 | 108.1 | 108 | 80-120 | 1 | 0-20 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



Quality Control - Spike/Spike Duplicate

| Environmental Science International, Inc. | Date Received: | 07/26/13 |
|---|----------------|--------------------|
| 354 Uluniu Street, Suite 304 | Work Order: | 13-07-1752 |
| Kailua, HI 96734-2500 | Preparation: | EPA 3510C |
| | Method: | EPA 8270C SIM PAHs |
| Project: Red Hill LTM 112066 | | Page 3 of 5 |

| Quality Control Sample ID | | Matrix | | Instrument | Date P | repared | Date Analyzed | MS | /MSD Batch | Number |
|---------------------------|------------------------|-----------------------|------------------------------|---------------------------|--------------------------|---------------------|-----------------|------------|---------------|-------------------|
| ES 034 | | Aqueou | Aqueous GC/MS AAA 07/29/13 0 | | 08/01/13 15:40 130729S12 | | | | | |
| Parameter | <u>Sample</u> Conc. | <u>Spike</u> Added | <u>MS</u> Conc. | <u>MS</u> <u>%Rec.</u> | <u>MSD</u> Conc. | <u>MSD</u> %Rec. | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Naphthalene | ND | 2.000 | 1.278 | 64 | 1.624 | 81 | 21-133 | 24 | 0-25 | |
| 2-Methylnaphthalene | ND | 2.000 | 1.155 | 58 | 1.466 | 73 | 21-140 | 24 | 0-25 | |
| 1-Methylnaphthalene | ND | 2.000 | 1.060 | 53 | 1.341 | 67 | 20-140 | 23 | 0-25 | |
| Acenaphthylene | ND | 2.000 | 1.241 | 62 | 1.584 | 79 | 33-145 | 24 | 0-25 | |
| Acenaphthene | ND | 2.000 | 1.242 | 62 | 1.585 | 79 | 49-121 | 24 | 0-25 | |
| Fluorene | ND | 2.000 | 1.298 | 65 | 1.680 | 84 | 59-121 | 26 | 0-25 | 4 |
| Phenanthrene | ND | 2.000 | 1.170 | 59 | 1.514 | 76 | 54-120 | 26 | 0-25 | 4 |
| Anthracene | ND | 2.000 | 1.049 | 52 | 1.276 | 64 | 27-133 | 20 | 0-25 | |
| Fluoranthene | ND | 2.000 | 1.165 | 58 | 1.475 | 74 | 26-137 | 24 | 0-25 | |
| Pyrene | ND | 2.000 | 1.264 | 63 | 1.622 | 81 | 18-168 | 25 | 0-25 | |
| Benzo (a) Anthracene | ND | 2.000 | 1.265 | 63 | 1.595 | 80 | 33-143 | 23 | 0-25 | |
| Chrysene | ND | 2.000 | 1.170 | 58 | 1.483 | 74 | 17-168 | 24 | 0-25 | |
| Benzo (k) Fluoranthene | ND | 2.000 | 1.229 | 61 | 1.632 | 82 | 24-159 | 28 | 0-25 | 4 |
| Benzo (b) Fluoranthene | ND | 2.000 | 1.307 | 65 | 1.684 | 84 | 24-159 | 25 | 0-25 | |
| Benzo (a) Pyrene | ND | 2.000 | 1.409 | 70 | 1.775 | 89 | 17-163 | 23 | 0-25 | |
| Indeno (1,2,3-c,d) Pyrene | ND | 2.000 | 1.284 | 64 | 1.634 | 82 | 10-171 | 24 | 0-25 | |
| Dibenz (a,h) Anthracene | ND | 2.000 | 1.107 | 55 | 1.436 | 72 | 10-219 | 26 | 0-25 | 4 |
| Benzo (g,h,i) Perylene | ND | 2.000 | 1.031 | 52 | 1.317 | 66 | 10-227 | 24 | 0-25 | |



| Environmental Science International, Inc. | Date Received: | 07/26/13 |
|---|----------------|-------------------|
| 354 Uluniu Street, Suite 304 | Work Order: | 13-07-1752 |
| Kailua, HI 96734-2500 | Preparation: | EPA 5030C |
| | Method: | GC/MS / EPA 8260B |
| Project: Red Hill LTM 112066 | | Page 4 of 5 |

Project: Red Hill LTM 112066

| E8 0.00CHAROW <thcharow< th="">CHAROWCHAROWCHAR</thcharow<> | Quality Control Sample ID | | Matrix | | Instrument | Date Pr | epared | Date Analyzed | MS | /MSD Batch | Number |
|--|-----------------------------|-------|--------|--------------------|--------------------|---------|---------------------|----------------|-----|------------|------------|
| Acetone88.3660.00187.2198175.617440.14060.203BenzeneND50.0051.0710247.449580-12070.20BromolichomethaneND50.0047.199447.659075.1310.2050.20BromolichomethaneND50.0057.13114101.7712430.14580.205BromolemethaneND50.0060.9712265.9111430.16570.205Catron TetrachlorideND50.0063.3010745.559960.1580.205ChiorothanaND50.0053.5810761.6813340.12540.205ChiorothaneND50.0053.5810751.6810340.12540.205DibromochioromethaneND50.0053.5810751.6810340.12550.2051.2-DibromochaneND50.0047.9084.7179480.1250.2051.2-DibromochaneND50.0047.929645.7270.13580.201.2-DibromochaneND50.0047.929645.729170.13580.201.2-DibromochaneND50.0047.929645.729170.13580.201.2-DibromochaneND50.00< | ES 034 | | Aqueou | us | GC/MS LL | 07/30/1 | 3 | 07/31/13 03:14 | 130 | 730S02 | |
| BarzaneND50.0047.1947.449580-12070.20BromodichloromethaneND50.0047.199444.769075.12050.20BromodicmND50.0047.199444.769075.12050.20BromodicmND50.0060.9712255.9111430.15070.202-ButanoneND50.0060.9712255.9111465.14070.20Carbon TetrachlorideND50.0048.879444.759080-13570.20ChloroberzaneND50.0053.3010749.559960-13570.20ChlorobertaneND50.0053.9110047.909660-13550.20ChlorobertaneND50.0050.1910047.909660-13550.201/2-Dibromo-3-ChloropropaneND50.0049.719147.199480.12050.201/2-Dibromo-3-ChloropropaneND50.0048.909247.378770.12050.201/2-Dibromo-3-ChloropropaneND50.0048.909345.799270.13580.201/2-DibroberzeneND50.0045.9110348.907010.2011/2-DibroberzeneND50.0051.5110348.907010.201 <td< th=""><th>Parameter</th><th></th><th></th><th><u>MS</u> Conc.</th><th><u>MS</u> %Rec.</th><th></th><th><u>MSD</u> %Rec.</th><th>%Rec. CL</th><th>RPD</th><th>RPD CL</th><th>Qualifiers</th></td<> | Parameter | | | <u>MS</u> Conc. | <u>MS</u> %Rec. | | <u>MSD</u> %Rec. | %Rec. CL | RPD | RPD CL | Qualifiers |
| BromodichloromethaneND50.0047.199444.769075-12050-20BromoderhaneND50.0045.199647.659571-13010-202-ButanoneND50.0060.9712256.9111430-15070-20Carbon TetrachlorideND50.0048.879845.619165-14070-20ChloroberaneND50.0064.879845.619166-13570-20ChloroberaneND50.0063.3010749.559366-13580-20ChloroberaneND50.0053.8110146.589365-13680-20ChloroberaneND50.0041.206239.6660-13550-2012-DichoroberaneND50.0047.199843.008676-12050-2012-DichoroberaneND50.0047.928947.7194600-20112-DichoroberaneND50.0047.928645.799270-13070-2012-DichoroberaneND50.0047.929445.799270-13070-2013-DichoroberaneND50.0051.5110347.459570-13070-2014-DichoroberaneND50.0051.5110348.819470-13070-20 <td< td=""><td>Acetone</td><td>88.36</td><td>50.00</td><td>187.2</td><td>198</td><td>175.6</td><td>174</td><td>40-140</td><td>6</td><td>0-20</td><td>3</td></td<> | Acetone | 88.36 | 50.00 | 187.2 | 198 | 175.6 | 174 | 40-140 | 6 | 0-20 | 3 |
| Bromotorm BromotertaneND50.0048.199647.659570-13010.20BromomethaneND50.0067.1311461.7712430-14580.202-ButanoneND50.0048.959444.759080-12050.20ChlorobenzeneND50.0048.879444.759080-12050.20ChlorobenzeneND50.0053.8010749.559960-13570.20ChlorobentaneND50.0053.8110751.6810340-12540.20ChlorobentaneND50.0053.8110751.6810340-12540.20L'Dohrom-3-ChloropropaneND50.0041.208239.688060.13030.201.2-DibromoethaneND50.0041.208243.738770-12050.201.2-DibromoethaneND50.0047.929645.429175-12540.201.2-DibromoethaneND50.0047.929645.429175-12550.201.2-DibromoethaneND50.0051.5110347.459570-13080.201.2-DibriorobenzeneND50.0051.5110347.819171-12550.201.2-DibriorobenzeneND50.0051.5110347.819171- | Benzene | ND | 50.00 | 51.07 | 102 | 47.44 | 95 | 80-120 | 7 | 0-20 | |
| BromomethaneND50.0057.1311461.7712430-14580-202-ButanoneND50.0060.9712256.9111430-16070-20Carbon TetrachiorideND50.0048.959445.619165.10070-20ChioroberzeneND50.0053.3010749.559960-13570-20ChiorobertaneND50.0053.3010746.569365.13580-20ChiorobertaneND50.0053.5810751.6810340-12540-20DibromochloromethaneND50.0045.0947.099660-13550-201.2-DibromochloromethaneND50.0041.208239.668051-13030-201.2-DibrioroberzeneND50.0041.208247.179480-12050-201.2-DibrioroberzeneND50.0047.829645.429175-12650-201.3-DibrioroberzeneND50.0051.5110347.459570-13080-201.2-DibrioroberzeneND50.0051.5110347.459570-13080-201.2-DibrioroberzeneND50.0051.5110346.849476-12560-201.2-DibrioropropaneND50.0051.5110346.849476- | Bromodichloromethane | ND | 50.00 | 47.19 | 94 | 44.76 | 90 | 75-120 | 5 | 0-20 | |
| 2-ButanoneND50.0060.9712256.9111430.15070.20Carbon TetrachlorideND50.0046.879445.759060.12050.20ChlorobharzeneND50.0046.879447.579060.12050.20ChlorothaneND50.0053.3010749.559860.13550.20ChlorothaneND50.0050.4310145.889366.13550.20DibromochloromethaneND50.0041.208239.868050.13030.201,2-DibromochaneND50.0041.208239.868050.13030.201,2-DibromochaneND50.0044.809043.008675.12550.201,3-DichlorobenzeneND50.0047.929645.429175.12550.201,4-DichlorobenzeneND50.0047.929645.429175.12550.201,4-DichlorobenzeneND50.0045.6110347.459570.13080.201,2-DichlorobenzeneND50.0051.5110347.459570.13080.201,2-DichlorobenzeneND50.0051.5110346.839470.12560.201,2-DichlorobenzeneND50.0051.5110346.849470. | Bromoform | ND | 50.00 | 48.19 | 96 | 47.65 | 95 | 70-130 | 1 | 0-20 | |
| Carbon TetrachlorideND50.0048.959845.619165.14070.20ChlorobenzeneND50.0068.879444.759060.13550.20ChlorobenzeneND50.0050.4310749.559960.13570.20ChlorobentaneND50.0050.4310145.889365.13580.20ChloromethaneND50.0051.5810751.6810340.12540.20L'2-Dibromo-3-ChloropropaneND50.0041.208239.668050.13030.201,2-DibromethaneND50.0044.009243.738770.12050.201,2-DibromethaneND50.0044.809043.008075.12550.201,3-DichlorobenzeneND50.0047.929645.429175.12550.201,4-DichlorobenzeneND50.0047.929645.429175.12580.201,1-DichloroethaneND50.0051.5110347.459470.13580.201,2-DichloroethaneND50.0051.5110348.309761.4070.201,2-DichloroethaneND50.0051.5110348.309761.4070.201,2-DichloroethaneND50.0051.5110348.3097 | Bromomethane | ND | 50.00 | 57.13 | 114 | 61.77 | 124 | 30-145 | 8 | 0-20 | |
| ChlorobenzeneND50.0046.879444.759080-12050.20ChlorotentaneND50.0053.3010749.559960-13570.20ChlorotentaneND50.0053.8010751.6810340-12540.20DibromochloromethaneND50.0053.5810751.6810340-12550.201,2-Dibromo-3-ChloropropaneND50.0041.208239.968050-13030.201,2-Dibromo-s-ChloropropaneND50.0040.719443.738770-12050.201,2-Dibromo-s-ChloropropaneND50.0044.889043.008675-12540.201,3-DichlorobenzeneND50.0047.929645.429175-12550.201,1-DichloroethaneND50.0047.929645.429170-13580.201,1-DichloroethaneND50.0051.5110347.459570-13080.201,1-DichloroethaneND50.0051.5110347.459570-13080.201,2-DichloroptaneND50.0051.5110347.459570-13080.201,2-DichloroptaneND50.0051.5110348.3091.0151.4040.201,2-DichloroptaneND50.0051.55103< | 2-Butanone | ND | 50.00 | 60.97 | 122 | 56.91 | 114 | 30-150 | 7 | 0-20 | |
| ChloroethaneND50.0053.3010749.559960-13570.20ChloroofromND50.0050.4310146.589365-13580.20ChloroothaneND50.0053.5810751.6810340.12540.20DibromochloromethaneND50.0041.208239.968050-13030.201,2-Dibromod-3-ChloropropaneND50.0049.719947.179480-12050.201,2-DibrombehaneND50.0044.809043.008675-12550.201,3-DichlorobenzeneND50.0044.809945.799270-13580.201,4-DichlorobenzeneND50.0049.609945.799270-13580.201,1-DichloroethaneND50.0051.5110347.509570-10380.20-1,2-DichloroethaneND50.0051.5110348.309760-13070.20-1,2-DichloroethaneND50.0051.6810348.309760-14070.20-1,2-DichloroetheneND50.0051.5110348.819870-13080.20-1,2-DichloroetheneND50.0051.5610348.819870-13080.20-1,2-DichloroetheneND50.0051.5510348.81 <td>Carbon Tetrachloride</td> <td>ND</td> <td>50.00</td> <td>48.95</td> <td>98</td> <td>45.61</td> <td>91</td> <td>65-140</td> <td>7</td> <td>0-20</td> <td></td> | Carbon Tetrachloride | ND | 50.00 | 48.95 | 98 | 45.61 | 91 | 65-140 | 7 | 0-20 | |
| ChloroformND50.0050.4310146.589365-13580-20ChloromethaneND50.0053.5810751.6810340-12540-20DibromochloromethaneND50.0050.1910047.909660-13550-201,2-Dibromo-3-ChloropropaneND50.0041.208239.968050-13030-201,2-Dibromo-thaneND50.0041.209243.738770-12050-201,2-DichlorobenzeneND50.0044.809043.008675-12550-201,3-DichlorobenzeneND50.0047.929645.709270-13580-201,1-DichloroethaneND50.0047.929645.709270-13580-201,1-DichloroethaneND50.0051.5110347.459570-13070-201,2-DichloroethaneND50.0051.5110348.309760.4070-201,2-DichloroetheneND50.0051.5510348.319870-13070-201,2-DichloroptheneND50.0051.5510348.319870-13070-201,2-DichloroptheneND50.0051.5510348.319870-13050-201,2-DichloroptheneND50.0051.5510348.31 <td>Chlorobenzene</td> <td>ND</td> <td>50.00</td> <td>46.87</td> <td>94</td> <td>44.75</td> <td>90</td> <td>80-120</td> <td>5</td> <td>0-20</td> <td></td> | Chlorobenzene | ND | 50.00 | 46.87 | 94 | 44.75 | 90 | 80-120 | 5 | 0-20 | |
| ChloromethaneND50.0053.5810751.6810340-12540-20DibromochloromethaneND50.0050.1910047.909660-13550-201,2-Dibromo-3-ChloropropaneND50.0041.208239.968050-13030-201,2-Dibromo-3-ChloropropaneND50.0049.719947.179480-12050-201,2-DibrlorobenzeneND50.0049.709243.738775-12550-201,3-DichlorobenzeneND50.0047.929645.429175-12550-201,4-DichlorobenzeneND50.0049.609945.799270-13080-201,1-DichloroethaneND50.0051.5110347.929470-13080-20-1,2-DichloroethaneND50.0051.5110347.929470-13080-20-1,2-DichloroethaneND50.0051.5110348.309760-14070-20-1,2-DichloroethaneND50.0051.5110348.309760-14070-20-1,2-DichloroethaneND50.0051.5110348.309760-14070-20-1,2-DichloroethaneND50.0051.5510348.309475-12580-20-1,2-DichloroethaneND50.0051.55 <td>Chloroethane</td> <td>ND</td> <td>50.00</td> <td>53.30</td> <td>107</td> <td>49.55</td> <td>99</td> <td>60-135</td> <td>7</td> <td>0-20</td> <td></td> | Chloroethane | ND | 50.00 | 53.30 | 107 | 49.55 | 99 | 60-135 | 7 | 0-20 | |
| DibromochloromethaneND50.0050.1910047.909660-13550-201,2-Dibromo-3-ChloropropaneND50.0041.208239.968050-13030-201,2-DibromothaneND50.0049.719947.179480-12050-201,2-DichlorobenzeneND50.0046.009243.738770-12050-201,3-DichlorobenzeneND50.0047.929645.429175-12550-201,4-DichlorobenzeneND50.0049.609945.799270-13080-201,1-DichloroethaneND50.0051.5110347.699470-13070-201,1-DichloroethaneND50.0050.4310146.949470-13070-20-1,2-DichloroethaneND50.0050.7110146.849470-13070-20-1,2-DichloroethaneND50.0050.7110146.849475-12580-20-1,3-DichloropropaneND50.0051.5510348.819870-13070-20-1,3-DichloropropaneND50.0052.4110150.3310155-14040-20-1,3-DichloropropaneND50.0052.4110550.3310155-14040-20-1,3-DichloropropaneND50.0052.41 <td>Chloroform</td> <td>ND</td> <td>50.00</td> <td>50.43</td> <td>101</td> <td>46.58</td> <td>93</td> <td>65-135</td> <td>8</td> <td>0-20</td> <td></td> | Chloroform | ND | 50.00 | 50.43 | 101 | 46.58 | 93 | 65-135 | 8 | 0-20 | |
| 1,2-Dibromo-3-ChloropropaneND50.0041.208239.968050-13030-201,2-DibromoethaneND50.0046.009243.738770-12050-201,3-DichlorobenzeneND50.0044.889043.008675-12540-201,4-DichlorobenzeneND50.0047.929645.429175-12550-201,4-DichlorobenzeneND50.0049.609045.799270-13080-201,2-DichlorobetnaneND50.0049.609447.459570-13080-201,2-DichloroethaneND50.0051.5110347.459470-13070-201,2-DichloroethaneND50.0051.6110346.949470-13070-201,2-DichloroethaneND50.0051.6810348.819870-13080-20t-1,2-DichloroethaneND50.0051.5510348.119870-13070-20t-1,2-DichloropropaneND50.0051.5510348.119870-13080-20t-1,3-DichloropropaneND50.0051.5510348.119870-13080-20t-1,3-DichloropropeneND50.0051.5510348.119265-13070-20t-1,3-DichloropropeneND50.0052.54 <td>Chloromethane</td> <td>ND</td> <td>50.00</td> <td>53.58</td> <td>107</td> <td>51.68</td> <td>103</td> <td>40-125</td> <td>4</td> <td>0-20</td> <td></td> | Chloromethane | ND | 50.00 | 53.58 | 107 | 51.68 | 103 | 40-125 | 4 | 0-20 | |
| 1.2-DibromoethaneND50.0049.719947.179480-12050-201.2-DichlorobenzeneND50.0046.009243.738770-12050-201.3-DichlorobenzeneND50.0044.889043.008675-12540-201.4-DichlorobenzeneND50.0047.929645.42917512550-201.1-DichloroethaneND50.0049.609945.799270-13580-201.1-DichloroethaneND50.0050.4310146.929470-13070-201.1-DichloroethaneND50.0050.4310146.929470-13070-201.1-DichloroetheneND50.0051.6810348.309760-14070-201.2-DichloroetheneND50.0051.5110348.309475-12580-201.2-DichloroptopaneND50.0051.5110348.319876-13050-201.3-DichloroptopeneND50.0040.488138.877855-14040-201.1-DichloroptopeneND50.0047.229446.139260-13520-201.1-DichloroptopeneND50.0045.5411050.3310155-14090-201.1-DichloroptopeneND50.0052.5410549 | Dibromochloromethane | ND | 50.00 | 50.19 | 100 | 47.90 | 96 | 60-135 | 5 | 0-20 | |
| 1.2.DichlorobenzeneND50.0046.009243.738770.12050.201.3.DichlorobenzeneND50.0044.889043.008675.12540.201.4.DichlorobenzeneND50.0047.929645.429175.12550.201.1.DichloroethaneND50.0049.609945.799270.13580.201.2.DichloroethaneND50.0051.5110347.459570.13070.201.1.DichloroethaneND50.0049.9910146.929470.13070.20t.1.2.DichloroethaneND50.0049.9910046.949470.12060.20t.1.2.DichloroetheneND50.0050.7110146.849475.12580.20t.1.3.DichloropropaneND50.0051.5110348.819870.13050.20t.1.3.DichloropropeneND50.0051.5510348.819870.13050.20t.1.3.DichloropropeneND50.0052.4110150.3310155.14040.20t.1.3.DichloropropeneND50.0052.4110349.319865.13570.20t.1.4.DichloroethaneND50.0052.5110349.139865.13570.20t.1.3.DichloropropeneND50.0052.51 </td <td>1,2-Dibromo-3-Chloropropane</td> <td>ND</td> <td>50.00</td> <td>41.20</td> <td>82</td> <td>39.96</td> <td>80</td> <td>50-130</td> <td>3</td> <td>0-20</td> <td></td> | 1,2-Dibromo-3-Chloropropane | ND | 50.00 | 41.20 | 82 | 39.96 | 80 | 50-130 | 3 | 0-20 | |
| 1.3-DichlorobenzeneND50.0044.889043.008675-12540-201.4-DichlorobenzeneND50.0047.929645.429175-12550-201.1-DichloroethaneND50.0049.609945.799270-13580-201.2-DichloroethaneND50.0051.5110347.459570-13080-201.1-DichloroethaneND50.0050.4310146.929470-13070-20c-1.2-DichloroetheneND50.0051.6810348.309760-14070-20t-1.2-DichloroetheneND50.0051.5510348.819870-13050-20t-1.2-DichloroptopaneND50.0051.5510348.819475-12580-20c-1.3-DichloroptopeneND50.0051.5510348.819870-13050-20t-1.3-DichloroptopeneND50.0040.488138.877855-14040-20t-1.3-DichloroptopeneND50.0047.029446.119260-13520-20t-1.3-DichloroptopeneND50.0055.2411050.3310155-14040-20t-1.3-DichloroethaneND50.0055.2410549.159865-13570-20t-1.3-DichloroptopeneND50.0051.35< | 1,2-Dibromoethane | ND | 50.00 | 49.71 | 99 | 47.17 | 94 | 80-120 | 5 | 0-20 | |
| 1,4-DichlorobenzeneND50.0047.929645.429175-12550-201,1-DichloroethaneND50.0049.609945.799270-13580-201,2-DichloroethaneND50.0051.5110347.459570-13080-201,1-DichloroethaneND50.0050.4310146.929470-13070-20c-1,2-DichloroetheneND50.0050.4310146.929470-12560-20t-1,2-DichloroetheneND50.0050.7110348.309760-14070-20t-1,2-DichloroptopaneND50.0050.7110146.849470-13280-20t-1,3-DichloroptopeneND50.0051.5510348.819870-13050-20t-1,3-DichloroptopeneND50.0051.5510348.819870-13050-20t-1,3-DichloroptopeneND50.0040.488138.877855-14040-20t-1,3-DichloroptopeneND50.0047.029446.119260-13520-20Methylene ChlorideND50.0052.5110349.119561-13570-201,1,1,2-TetrachloroethaneND50.0051.3510349.119186.13570-201,1,1,2-TetrachloroethaneND50.00 <t< td=""><td>1,2-Dichlorobenzene</td><td>ND</td><td>50.00</td><td>46.00</td><td>92</td><td>43.73</td><td>87</td><td>70-120</td><td>5</td><td>0-20</td><td></td></t<> | 1,2-Dichlorobenzene | ND | 50.00 | 46.00 | 92 | 43.73 | 87 | 70-120 | 5 | 0-20 | |
| 1,1-DichloroethaneND50.0049.609945.799270-13580-201,2-DichloroethaneND50.0051.5110347.459570-13080-201,1-DichloroethaneND50.0050.4310146.929470-13070-20c-1,2-DichloroethaneND50.0049.9910046.949470-12560-20t-1,2-DichloropthaneND50.0051.6810348.309760-14070-201,2-DichloroptpaneND50.0051.7510348.819870-13050-20c-1,3-DichloroptpaneND50.0051.5510348.819870-13050-20t-1,3-DichloroptpaneND50.0049.488138.877855-14040-20t-1,3-DichloroptpaneND50.0049.089846.939475-15240-20t-1,3-DichloroptpaneND50.0049.089846.939475-15240-20t-1,3-DichloroptpaneND50.0049.089846.939475-15240-20t-1,3-DichloroptpaneND50.0052.4411050.3310155-14090-20t-1,3-DichloropthaneND50.0052.4411050.3310155-14040-20t-1,1,1-TettachloroethaneND50.0051.55< | 1,3-Dichlorobenzene | ND | 50.00 | 44.88 | 90 | 43.00 | 86 | 75-125 | 4 | 0-20 | |
| 1,2-DichloroethaneND50.0051.5110347.459570-13080-201,1-DichloroetheneND50.0050.4310146.929470-13070-20c-1,2-DichloroetheneND50.0049.9910046.949470-12560-20t-1,2-DichloroetheneND50.0051.6810348.309760-14070-201,2-DichloroptopaneND50.0051.5510348.819870-13050-20c-1,3-DichloroptopeneND50.0051.5510348.819870-13050-20t-1,3-DichloroptopeneND50.0051.5510348.819875-12540-20t-1,3-DichloroptopeneND50.0049.089846.939475-12540-20t-1,3-DichloroptopeneND50.0055.2411050.3310155.14090-20t-1,3-DichloroptopeneND50.0055.2411050.3310155.14090-20t-1,1,2-PertachloroethaneND50.0051.5510549.159866.13570-201,1,2-TetrachloroethaneND50.0051.5510549.159865.136640-203,41,1,2-TetrachloroethaneND50.0051.351060.695.0165.13570-201,1,2-Tetrachloroethane | 1,4-Dichlorobenzene | ND | 50.00 | 47.92 | 96 | 45.42 | 91 | 75-125 | 5 | 0-20 | |
| 1,1-DichloroetheneND50.0050.4310146.929470-13070-20c-1,2-DichloroetheneND50.0049.9910046.949470-1256020t-1,2-DichloroetheneND50.0051.6810348.309760-14070201,2-DichloroptopaneND50.0050.7110146.849475-1258020c-1,3-DichloroptopeneND50.0051.5510348.819870-1305020t-1,3-DichloroptopeneND50.0040.488138.877855-1404020EthylbenzeneND50.0049.089846.939475-1254020Methyl-2-PentanoneND50.0047.029446.119260-13520201,1,2-TetrachloroethaneND50.0052.5110549.159865-13570201,1,2-TetrachloroethaneND50.0051.3510349.319980-13040201,1,2-TetrachloroethaneND50.0076.4515372.0014445-150602031,1,2-TetrachloroethaneND50.0076.4515372.0014445-150602031,2,4-TrichloroethaneND50.0048.199645.869265-13550201,2,4-TrichloroethaneND50.00 | 1,1-Dichloroethane | ND | 50.00 | 49.60 | 99 | 45.79 | 92 | 70-135 | 8 | 0-20 | |
| c.1,2-DichloroetheneND50.0049.9910046.949470-12560-20t-1,2-DichloroetheneND50.0051.6810348.309760-14070-201,2-DichloropropaneND50.0050.7110146.849475-12580-20c-1,3-DichloropropeneND50.0051.5510348.819870-13050-20t-1,3-DichloropropeneND50.0040.488138.877855-14040-20t-1,3-DichloropropeneND50.0049.089846.939475-12540-20EthylbenzeneND50.0047.029446.119260-13520-204-Methyl-2-PentanoneND50.0047.029446.119260-13520-205,yreneND50.0052.5110549.159865-13570-201,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-203,41,1,2-ZetrachloroethaneND50.0076.4515372.2014445-15060-203,41,2,4-TrichloroethaneND50.0076.4515372.2014445-15060-203,41,2,4-TrichloroethaneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND | 1,2-Dichloroethane | ND | 50.00 | 51.51 | 103 | 47.45 | 95 | 70-130 | 8 | 0-20 | |
| t1.2-DichloroetheneND50.0051.6810348.309760-14070-201,2-DichloropropaneND50.0050.7110146.849475-12580-20c-1,3-DichloropropeneND50.0051.5510348.819870-13050-20t-1,3-DichloropropeneND50.0040.488138.877855-14040-20t-1,3-DichloropropeneND50.0049.089846.939475-12540-20EthylbenzeneND50.0055.2411050.3310155-14090-204-Methyl-2-PentanoneND50.0047.029446.119260-13520-20StyreneND50.0052.5110349.319980-13040-201,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2-TetrachloroethaneND50.0076.4515372.2014445-15060-20341,1,2-TetrachloroethaneND50.0048.189945.899275-12070-201,2,4-TrichlorobenzeneND50.0048.199645.869265-13550-201,2,4-TrichloroethaneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND50.00 | 1,1-Dichloroethene | ND | 50.00 | 50.43 | 101 | 46.92 | 94 | 70-130 | 7 | 0-20 | |
| 1,2-DichloropropaneND50.0050.7110146.849475-12580-20c-1,3-DichloropropeneND50.0051.5510348.819870-13050-20t-1,3-DichloropropeneND50.0040.488138.877855-14040-20EthylbenzeneND50.0049.089846.939475-12540-20Methylene ChlorideND50.0047.029460.3310155-14090-204-Methyl-2-PentanoneND50.0052.5110549.159865-13570-20StyreneND50.0051.3510349.159865-13570-201,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2-TetrachloroethaneND50.0051.3510349.319165-13060-203.4TetrachloroethaneND50.0076.4515372.2014445-15060-203.4TolueneND50.0048.199645.869255-13070-201,2,4-TrichlorobehaneND50.0048.199644.879065-13070-201,1,1-TrichloroethaneND50.0048.199644.879065-13070-201,1,1-TrichloroethaneND50.0048.19 <td>c-1,2-Dichloroethene</td> <td>ND</td> <td>50.00</td> <td>49.99</td> <td>100</td> <td>46.94</td> <td>94</td> <td>70-125</td> <td>6</td> <td>0-20</td> <td></td> | c-1,2-Dichloroethene | ND | 50.00 | 49.99 | 100 | 46.94 | 94 | 70-125 | 6 | 0-20 | |
| c-1,3-DichloropropeneND50.0051.5510348.819870-13050-20t-1,3-DichloropropeneND50.0040.488138.877855-14040-20EthylbenzeneND50.0049.089846.939475-12540-20Methylene ChlorideND50.0055.2411050.3310155-14090-204-Methyl-2-PentanoneND50.0055.2411050.3310155-13520-20StyreneND50.0052.5110549.159865-13570-201,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2,2-TetrachloroethaneND50.0051.3510349.319980-130640-203,4TetrachloroethaneND50.0076.4515372.2014445-150640-203,4TolueneND50.0049.389945.899275-12070-201,1,2,4-TrichlorobenzeneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND50.0048.199644.879065-13070-201,1,1-TrichloroethaneND50.0048.019644.879065-13050-201,1,1-TrichloroethaneND50.0048.01 <td>t-1,2-Dichloroethene</td> <td>ND</td> <td>50.00</td> <td>51.68</td> <td>103</td> <td>48.30</td> <td>97</td> <td>60-140</td> <td>7</td> <td>0-20</td> <td></td> | t-1,2-Dichloroethene | ND | 50.00 | 51.68 | 103 | 48.30 | 97 | 60-140 | 7 | 0-20 | |
| t 1,3-DichloropropeneND50.0040.488138.877855-14040-20EthylbenzeneND50.0049.089846.939475-12540-20Methylene ChlorideND50.0055.2411050.3310155-14090-204-Methyl-2-PentanoneND50.0047.029446.119260-13520-20StyreneND50.0052.5110549.159865-13570-201,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2,2-TetrachloroethaneND50.0076.4515372.2014445-15060-203,4TolueneND50.0048.199645.899275-12070-201,1,1,1-TrichloroethaneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-201,1,1-TrichloroethaneND50.0048.019644.869050-14050-201,1,1-TrichloroethaneND50.0048.0196 </td <td>1,2-Dichloropropane</td> <td>ND</td> <td>50.00</td> <td>50.71</td> <td>101</td> <td>46.84</td> <td>94</td> <td>75-125</td> <td>8</td> <td>0-20</td> <td></td> | 1,2-Dichloropropane | ND | 50.00 | 50.71 | 101 | 46.84 | 94 | 75-125 | 8 | 0-20 | |
| EthylbenzeneND50.0049.089846.939475-12540-20Methylene ChlorideND50.0055.2411050.3310155-14090-204-Methyl-2-PentanoneND50.0047.029446.119260-13520-20StyreneND50.0052.5110549.159865-13570-201,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2,2-TetrachloroethaneND50.000.358510.6950165-130640-203,4TetrachloroethaneND50.0076.4515372.2014445-15060-203TolueneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND50.0048.199644.879065-13070-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-201,1,1-TrichloroethaneND50.0047.459544.969050-14050-20Hexachloro-1,3-ButadieneND50.0047.459544.969050-14050-20 | c-1,3-Dichloropropene | ND | 50.00 | 51.55 | 103 | 48.81 | 98 | 70-130 | 5 | 0-20 | |
| Methylene ChlorideND50.0055.2411050.3310155-14090-204-Methyl-2-PentanoneND50.0047.029446.119260-13520-20StyreneND50.0052.5110549.159865-13570-201,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2,2-TetrachloroethaneND50.000.358510.6950165-130640-203,4TetrachloroethaneND50.0076.4515372.2014445-15060-203TolueneND50.0049.389945.899275-12070-201,2,4-TrichloroethaneND50.0048.199644.879065-13050-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-201,1,1-TrichloroethaneND50.0047.459544.969050-14050-20 | t-1,3-Dichloropropene | ND | 50.00 | 40.48 | 81 | 38.87 | 78 | 55-140 | 4 | 0-20 | |
| 4-Methyl-2-PentanoneND50.0047.029446.119260-13520-20StyreneND50.0052.5110549.159865-13570-201,1,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2,2-TetrachloroethaneND50.000.358510.6950165-130640-203,4TetrachloroethaneND50.0076.4515372.2014445-15060-203TolueneND50.0049.389945.899275-12070-201,2,4-TrichlorobenzeneND50.0048.199644.879065-13550-201,1,1-TrichloroethaneND50.0047.459544.969050-14050-20 | Ethylbenzene | ND | 50.00 | 49.08 | 98 | 46.93 | 94 | 75-125 | 4 | 0-20 | |
| StyreneND50.0052.5110549.159865-13570-201,1,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2,2-TetrachloroethaneND50.000.358510.6950165-130640-203,4TetrachloroethaneND50.0076.4515372.2014445-15060-203TolueneND50.0049.389945.899275-12070-201,2,4-TrichlorobenzeneND50.0048.199644.879065-13550-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-20Hexachloro-1,3-ButadieneND50.0047.459544.969050-14050-20 | Methylene Chloride | ND | 50.00 | 55.24 | 110 | 50.33 | 101 | 55-140 | 9 | 0-20 | |
| 1,1,2-TetrachloroethaneND50.0051.3510349.319980-13040-201,1,2,2-TetrachloroethaneND50.000.358510.6950165-130640-203,4TetrachloroetheneND50.0076.4515372.2014445-15060-203TolueneND50.0049.389945.899275-12070-201,2,4-TrichloroethaneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-20Hexachloro-1,3-ButadieneND50.0047.459544.969050-14050-20 | 4-Methyl-2-Pentanone | ND | 50.00 | 47.02 | 94 | 46.11 | 92 | 60-135 | 2 | 0-20 | |
| 1,1,2,2-TetrachloroethaneND50.000.358510.6950165-130640-203,4TetrachloroetheneND50.0076.4515372.2014445-15060-203TolueneND50.0049.389945.899275-12070-201,2,4-TrichlorobenzeneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-20Hexachloro-1,3-ButadieneND50.0047.459544.969050-14050-20 | Styrene | ND | 50.00 | 52.51 | 105 | 49.15 | 98 | 65-135 | 7 | 0-20 | |
| TetrachloroetheneND50.0076.4515372.2014445-15060-203TolueneND50.0049.389945.899275-12070-201,2,4-TrichlorobenzeneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-20Hexachloro-1,3-ButadieneND50.0047.459544.969050-14050-20 | 1,1,1,2-Tetrachloroethane | ND | 50.00 | 51.35 | 103 | 49.31 | 99 | 80-130 | 4 | 0-20 | |
| TolueneND50.0049.389945.899275-12070-201,2,4-TrichlorobenzeneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-20Hexachloro-1,3-ButadieneND50.0047.459544.969050-14050-20 | 1,1,2,2-Tetrachloroethane | ND | 50.00 | 0.3585 | 5 1 | 0.6950 | 1 | 65-130 | 64 | 0-20 | 3,4 |
| 1,2,4-TrichlorobenzeneND50.0048.199645.869265-13550-201,1,1-TrichloroethaneND50.0048.019644.879065-13070-20Hexachloro-1,3-ButadieneND50.0047.459544.969050-14050-20 | Tetrachloroethene | ND | 50.00 | 76.45 | 153 | 72.20 | 144 | 45-150 | 6 | 0-20 | 3 |
| 1,1,1-TrichloroethaneND50.0048.019644.879065-13070-20Hexachloro-1,3-ButadieneND50.0047.459544.969050-14050-20 | Toluene | ND | 50.00 | 49.38 | 99 | 45.89 | 92 | 75-120 | 7 | 0-20 | |
| Hexachloro-1,3-Butadiene ND 50.00 47.45 95 44.96 90 50-140 5 0-20 | 1,2,4-Trichlorobenzene | ND | 50.00 | 48.19 | 96 | 45.86 | 92 | 65-135 | 5 | 0-20 | |
| | 1,1,1-Trichloroethane | ND | 50.00 | 48.01 | 96 | 44.87 | 90 | 65-130 | 7 | 0-20 | |
| 1,1,2-Trichloroethane ND 50.00 50.27 101 47.95 96 75-125 5 0-20 | Hexachloro-1,3-Butadiene | ND | 50.00 | 47.45 | 95 | 44.96 | 90 | 50-140 | 5 | 0-20 | |
| | 1,1,2-Trichloroethane | ND | 50.00 | 50.27 | 101 | 47.95 | 96 | 75-125 | 5 | 0-20 | |

RPD: Relative Percent Difference. CL: Control Limits



| Environmental Science Internat | tional, Inc. | | | Date F | Received: | | | | | 07/26/13 |
|--------------------------------|------------------------|-----------------------|--------------------|---------------------------|---------------------|---------------------|-----------------|------------|-----------|-------------------|
| 354 Uluniu Street, Suite 304 | | | | Work (| Order: | | | | 1 | 3-07-1752 |
| Kailua, HI 96734-2500 | | | | Prepa | ation: | | | | E | PA 5030C |
| | | | | Metho | d: | | | C | GC/MS / E | PA 8260B |
| Project: Red Hill LTM 112066 | | | | | | | | | Page | 5 of 5 |
| Parameter | <u>Sample</u> Conc. | <u>Spike</u> Added | <u>MS</u> Conc. | <u>MS</u> <u>%Rec.</u> | <u>MSD</u> Conc. | <u>MSD</u> %Rec. | <u>%Rec. CL</u> | <u>RPD</u> | RPD CL | <u>Qualifiers</u> |
| Trichloroethene | ND | 50.00 | 87.33 | 175 | 80.84 | 162 | 70-125 | 8 | 0-20 | 3 |
| 1,2,3-Trichloropropane | ND | 50.00 | 50.03 | 100 | 48.01 | 96 | 75-125 | 4 | 0-20 | |
| Vinyl Chloride | ND | 50.00 | 50.64 | 101 | 47.81 | 96 | 50-145 | 6 | 0-20 | |
| p/m-Xylene | ND | 100.0 | 98.53 | 99 | 93.79 | 94 | 75-130 | 5 | 0-20 | |
| o-Xylene | ND | 50.00 | 48.14 | 96 | 45.54 | 91 | 80-120 | 6 | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | ND | 50.00 | 42.97 | 86 | 41.53 | 83 | 65-125 | 3 | 0-20 | |

Return to Contents

RPD: Relative Percent Difference. CL: Control Limits



| Environmental Science International, Inc. | Date Received: | 07/26/13 |
|---|----------------|-----------------|
| 354 Uluniu Street, Suite 304 | Work Order: | 13-07-1752 |
| Kailua, HI 96734-2500 | Preparation: | EPA 3020A Total |
| | Method: | EPA 6020 |
| Project: Red Hill LTM 112066 | | Page 1 of 1 |

| Quality Control Sample ID | Matrix | Instrument | Date Prepa | red Date Anal | yzed PDS/P | DSD Batch Number |
|---------------------------|--------------|-------------|-------------|---------------|-----------------|-------------------|
| ES 034 | Aqueous | ICP/MS 04 | 07/31/13 00 | 0:00 07/31/13 | 17:43 130731 | S02 |
| Parameter | Sample Conc. | Spike Added | PDS Conc. | PDS %Rec. | <u>%Rec. CL</u> | <u>Qualifiers</u> |
| Lead | ND | 100.0 | 106.1 | 106 | 75-125 | |



RPD: Relative Percent Difference. CL: Control Limits



| Environmental Science International, Inc. | Date Received: | 07/26/13 |
|---|----------------|---------------|
| 354 Uluniu Street, Suite 304 | Work Order: | 13-07-1752 |
| Kailua, HI 96734-2500 | Preparation: | EPA 3510C |
| | Method: | EPA 8015B (M) |
| Project: Red Hill LTM 112066 | | Page 1 of 5 |

| Quality Control Sample ID | | Matrix | | Instrument | Date Prepa | red Date | Analyzed | LCS/LCSD Ba | atch Number |
|---------------------------|-----------------------|--------------|----------------------------|---------------|----------------------|-----------------|------------|---------------|-------------------|
| 099-15-516-53 | | Aqueou | S | GC 45 | 07/30/13 | 07/31 | /13 01:52 | 130730B09 | |
| Parameter | <u>Spike</u> Added | LCS Conc. | <u>LCS</u> <u>%Rec.</u> | LCSD Conc. | <u>LCSD</u> %Rec. | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| TPH as Diesel | 4000 | 4379 | 109 | 4127 | 103 | 60-132 | 6 | 0-11 | |

Project: Red Hill LTM 112066

| Quality Control Sample ID | Matrix | Instrument | Date Ana | lyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|-----------|--------------|------------------|
| 099-14-497-42 | Aqueous | ICP/MS 04 | 07/31/13 | 17:31 | 130731L02D |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | <u>%Rec.</u> | CL Qualifiers |
| Lead | 100.0 | 94.32 | 94 | 80-120 | 1 |

Method:

EPA 6020

Page 2 of 5

alscience nvironmental aboratories, Inc.

| Environmental Science International, Inc. | Date Received: | 07/26/13 |
|---|----------------|--------------------|
| | Date Received. | 07/20/13 |
| 354 Uluniu Street, Suite 304 | Work Order: | 13-07-1752 |
| Kailua, HI 96734-2500 | Preparation: | EPA 3510C |
| | Method: | EPA 8270C SIM PAHs |
| Project: Red Hill LTM 112066 | | Page 3 of 5 |

Quality Control - LCS

Project: Red Hill LTM 112066

| Quality Control Sample ID | Matrix | Instrument | Date Ana | alyzed | LCS Batch Number |
|---------------------------|-------------|-----------------|-----------|----------------|-----------------------------|
| 099-15-148-17 | Aqueous | GC/MS AAA | 08/01/13 | 15:14 | 130729L12 |
| Parameter | Spike Added | Conc. Recovered | LCS %Rec. | <u>%Rec. (</u> | <u>CL</u> <u>Qualifiers</u> |
| Naphthalene | 2.000 | 1.682 | 84 | 21-133 | |
| 2-Methylnaphthalene | 2.000 | 1.748 | 87 | 21-140 | |
| 1-Methylnaphthalene | 2.000 | 1.625 | 81 | 20-140 | |
| Acenaphthylene | 2.000 | 1.684 | 84 | 33-145 | |
| Acenaphthene | 2.000 | 1.736 | 87 | 55-121 | |
| Fluorene | 2.000 | 1.883 | 94 | 59-121 | |
| Phenanthrene | 2.000 | 1.878 | 94 | 54-120 | |
| Anthracene | 2.000 | 1.617 | 81 | 27-133 | |
| Fluoranthene | 2.000 | 1.900 | 95 | 26-137 | |
| Pyrene | 2.000 | 2.068 | 103 | 45-129 | |
| Benzo (a) Anthracene | 2.000 | 2.080 | 104 | 33-143 | |
| Chrysene | 2.000 | 1.998 | 100 | 17-168 | |
| Benzo (k) Fluoranthene | 2.000 | 2.317 | 116 | 24-159 | |
| Benzo (b) Fluoranthene | 2.000 | 2.263 | 113 | 24-159 | |
| Benzo (a) Pyrene | 2.000 | 2.295 | 115 | 17-163 | |
| Indeno (1,2,3-c,d) Pyrene | 2.000 | 2.148 | 107 | 25-175 | |
| Dibenz (a,h) Anthracene | 2.000 | 1.832 | 92 | 25-175 | |
| Benzo (g,h,i) Perylene | 2.000 | 1.723 | 86 | 25-157 | |

Total number of LCS compounds: 18 Total number of ME compounds: 0 Total number of ME compounds allowed: 1 LCS ME CL validation result: Pass

Return to Contents

| <i>nvironmental</i> | |
|---------------------|--|
| aboratories, Inc. | |

alscience

| Environmental Science International, Inc. | Date Received: | 07/26/13 |
|---|----------------|-------------------|
| 354 Uluniu Street, Suite 304 | Work Order: | 13-07-1752 |
| Kailua, HI 96734-2500 | Preparation: | EPA 5030C |
| | Method: | GC/MS / EPA 8260B |
| Project: Red Hill LTM 112066 | | Page 4 of 5 |

Project: Red Hill LTM 112066

| Quality Control Sample ID | | Matrix | | Instrument | Date Prepa | ared Date A | nalyzed | LCS/LCSD B | atch Number |
|-----------------------------|------------------------------|---------------------|----------------------------|---------------|-----------------------------|-----------------|------------|---------------|-------------------|
| 099-13-057-26 | | | IS | GC/MS LL | 07/30/13 | 07/30/ | 13 18:56 | 130730L02 | |
| Parameter | <u>Spike</u> <u>Added</u> | <u>LCS</u> Conc. | <u>LCS</u> <u>%Rec.</u> | LCSD Conc. | <u>LCSD</u> <u>%Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | <u>RPD CL</u> | <u>Qualifiers</u> |
| Acetone | 50.00 | 75.72 | 151 | N/A | N/A | 40-140 | N/A | 0-20 | |
| Benzene | 50.00 | 53.87 | 108 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Bromodichloromethane | 50.00 | 50.21 | 100 | N/A | N/A | 75-120 | N/A | 0-20 | |
| Bromoform | 50.00 | 58.22 | 116 | N/A | N/A | 70-130 | N/A | 0-20 | |
| Bromomethane | 50.00 | 58.56 | 117 | N/A | N/A | 30-145 | N/A | 0-20 | |
| 2-Butanone | 50.00 | 59.63 | 119 | N/A | N/A | 30-150 | N/A | 0-20 | |
| Carbon Tetrachloride | 50.00 | 50.51 | 101 | N/A | N/A | 65-140 | N/A | 0-20 | |
| Chlorobenzene | 50.00 | 50.85 | 102 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Chloroethane | 50.00 | 51.79 | 104 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Chloroform | 50.00 | 49.70 | 99 | N/A | N/A | 65-135 | N/A | 0-20 | |
| Chloromethane | 50.00 | 51.81 | 104 | N/A | N/A | 40-125 | N/A | 0-20 | |
| Dibromochloromethane | 50.00 | 54.93 | 110 | N/A | N/A | 60-135 | N/A | 0-20 | |
| 1,2-Dibromo-3-Chloropropane | 50.00 | 48.78 | 98 | N/A | N/A | 50-130 | N/A | 0-20 | |
| 1,2-Dibromoethane | 50.00 | 54.27 | 109 | N/A | N/A | 80-120 | N/A | 0-20 | |
| 1,2-Dichlorobenzene | 50.00 | 50.66 | 101 | N/A | N/A | 70-120 | N/A | 0-20 | |
| 1,3-Dichlorobenzene | 50.00 | 51.06 | 102 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,4-Dichlorobenzene | 50.00 | 53.41 | 107 | N/A | N/A | 75-125 | N/A | 0-20 | |
| 1,1-Dichloroethane | 50.00 | 50.23 | 100 | N/A | N/A | 70-135 | N/A | 0-20 | |
| 1,2-Dichloroethane | 50.00 | 49.86 | 100 | N/A | N/A | 70-130 | N/A | 0-20 | |
| 1,1-Dichloroethene | 50.00 | 49.81 | 100 | N/A | N/A | 70-130 | N/A | 0-20 | |
| c-1,2-Dichloroethene | 50.00 | 52.16 | 104 | N/A | N/A | 70-125 | N/A | 0-20 | |
| t-1,2-Dichloroethene | 50.00 | 53.49 | 107 | N/A | N/A | 60-140 | N/A | 0-20 | |
| 1,2-Dichloropropane | 50.00 | 53.34 | 107 | N/A | N/A | 75-125 | N/A | 0-20 | |
| c-1,3-Dichloropropene | 50.00 | 59.86 | 120 | N/A | N/A | 70-130 | N/A | 0-20 | |
| t-1,3-Dichloropropene | 50.00 | 46.32 | 93 | N/A | N/A | 55-140 | N/A | 0-20 | |
| Ethylbenzene | 50.00 | 54.54 | 109 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Methylene Chloride | 50.00 | 54.06 | 108 | N/A | N/A | 55-140 | N/A | 0-20 | |
| 4-Methyl-2-Pentanone | 50.00 | 54.61 | 109 | N/A | N/A | 60-135 | N/A | 0-20 | |
| Styrene | 50.00 | 55.76 | 112 | N/A | N/A | 65-135 | N/A | 0-20 | |
| 1,1,1,2-Tetrachloroethane | 50.00 | 53.71 | 107 | N/A | N/A | 80-130 | N/A | 0-20 | |
| 1,1,2,2-Tetrachloroethane | 50.00 | 54.37 | 109 | N/A | N/A | 65-130 | N/A | 0-20 | |
| Tetrachloroethene | 50.00 | 53.54 | 107 | N/A | N/A | 45-150 | N/A | 0-20 | |
| Toluene | 50.00 | 52.86 | 106 | N/A | N/A | 75-120 | N/A | 0-20 | |
| 1,2,4-Trichlorobenzene | 50.00 | 57.01 | 114 | N/A | N/A | 65-135 | N/A | 0-20 | |
| 1,1,1-Trichloroethane | 50.00 | 48.94 | 98 | N/A | N/A | 65-130 | N/A | 0-20 | |
| Hexachloro-1,3-Butadiene | 50.00 | 54.06 | 108 | N/A | N/A | 50-140 | N/A | 0-20 | |
| 1,1,2-Trichloroethane | 50.00 | 53.95 | 108 | N/A | N/A | 75-125 | N/A | 0-20 | |

RPD: Relative Percent Difference. CL: Control Limits



| Environmental Science Inter | rnational, In | C. | | Date Re | eceived: | | | | 07/26/13 |
|------------------------------|-----------------------|---------------------|----------------------------|---------------|----------------------|-----------------|------------|---------|-------------------|
| 354 Uluniu Street, Suite 304 | Work O | rder: | | 13-07-1752 | | | | | |
| Kailua, HI 96734-2500 | Prepara | ation: | | | EPA 5030C | | | | |
| | | | | Method | : | | | GC/MS / | EPA 8260B |
| Project: Red Hill LTM 11206 | 6 | | | | | | | Page | e 5 of 5 |
| Parameter | <u>Spike</u> Added | <u>LCS</u> Conc. | <u>LCS</u> <u>%Rec.</u> | LCSD Conc. | LCSD <u>%Rec.</u> | <u>%Rec. CL</u> | <u>RPD</u> | RPD CL | <u>Qualifiers</u> |
| Trichloroethene | 50.00 | 52.76 | 106 | N/A | N/A | 70-125 | N/A | 0-20 | |
| 1,2,3-Trichloropropane | 50.00 | 51.92 | 104 | N/A | N/A | 75-125 | N/A | 0-20 | |
| Vinyl Chloride | 50.00 | 50.66 | 101 | N/A | N/A | 50-145 | N/A | 0-20 | |
| p/m-Xylene | 100.0 | 106.1 | 106 | N/A | N/A | 75-130 | N/A | 0-20 | |
| o-Xylene | 50.00 | 51.18 | 102 | N/A | N/A | 80-120 | N/A | 0-20 | |
| Methyl-t-Butyl Ether (MTBE) | 50.00 | 47.32 | 95 | N/A | N/A | 65-125 | N/A | 0-20 | |
| Gasoline Range Organics | 1000 | 1013 | 101 | 939.9 | 94 | 80-120 | 8 | 0-20 | |

RPD: Relative Percent Difference. CL: Control Limits

Page 1 of 1



Work Order: 13-07-1752

| Method | Extraction | Chemist ID | Instrument | Analytical Location |
|--------------------|-----------------|------------|------------|---------------------|
| EPA 6020 | EPA 3020A Total | 598 | ICP/MS 04 | 1 |
| EPA 8015B (M) | EPA 3510C | 682 | GC 45 | 1 |
| EPA 8270C SIM PAHs | EPA 3510C | 773 | GC/MS AAA | 1 |
| GC/MS / EPA 8260B | EPA 5030C | 670 | GC/MS LL | 2 |

Location 1: 7440 Lincoln Way, Garden Grove, CA 92841 Location 2: 7445 Lampson Avenue, Garden Grove, CA 92841 alscience nvironmental aboratories, Inc.

Work Order: 13-07-1752

Page 1 of 1 Qualifiers Definition * See applicable analysis comment. Less than the indicated value. < Greater than the indicated value. > Surrogate compound recovery was out of control due to a required sample dilution. Therefore, the sample data was reported without further 1 clarification. 2 Surrogate compound recovery was out of control due to matrix interference. The associated method blank surrogate spike compound was in control and, therefore, the sample data was reported without further clarification. 3 Recovery of the Matrix Spike (MS) or Matrix Spike Duplicate (MSD) compound was out of control due to suspected matrix interference. The associated LCS recovery was in control. Δ The MS/MSD RPD was out of control due to suspected matrix interference. The PDS/PDSD or PES/PESD associated with this batch of samples was out of control due to suspected matrix interference. 5 6 Surrogate recovery below the acceptance limit. 7 Surrogate recovery above the acceptance limit. В Analyte was present in the associated method blank. ΒU Sample analyzed after holding time expired. ΒV Sample received after holding time expired. DL The Detection Limit (DL) is the smallest analyte concentration that can be demonstrated to be different from zero or a blank concentration at the 99% level of confidence. Е Concentration exceeds the calibration range. ET Sample was extracted past end of recommended max. holding time. HD The chromatographic pattern was inconsistent with the profile of the reference fuel standard. ICH Initial calibration verification recovery is above the control limit for this analyte. ICJ Initial calibration verification recovery is below the control limit for this analyte. IH Calibration verification recovery is above the control limit for this analyte. IJ Calibration verification recovery is below the control limit for this analyte. Analyte was detected at a concentration below the LOQ and above the DL. Reported value is estimated. J LOD The Limit of Detection (LOD) is the smallest amount or concentration of a substance that must be present in a sample in order to be detected at 99% confidence level. LOQ The Limit of Quantitation (LOQ) is the lowest concentration of a substance that produces a quantitative result within specified limits of precision and bias Q Spike recovery and RPD control limits do not apply resulting from the parameter concentration in the sample exceeding the spike concentration by a factor of four or greater. SG The sample extract was subjected to Silica Gel treatment prior to analysis. U Undetected at Detection Limit (DL) and is reported as less than the Limit of Detection (LOD). Х % Recovery and/or RPD out-of-range.

Glossary of Terms and Qualifiers

Ζ Analyte presence was not confirmed by second column or GC/MS analysis.

> Solid - Unless otherwise indicated, solid sample data is reported on a wet weight basis, not corrected for % moisture. All QC results are reported on a wet weight basis.

> Any parameter identified in 40CFR Part 136.3 Table II that is designated as "analyze immediately" with a holding time of <= 15 minutes (40CFR-136.3 Table II, footnote 4), is considered a "field" test and the reported results will be qualified as being received outside of the stated holding time unless received at the laboratory within 15 minutes of the collection time.

> A calculated total result (Example: Total Pesticides) is the summation of each component concentration and/or, if "J" flags are reported, estimated concentration. Component concentrations showing not detected (ND) are summed into the calculated total result as zero concentrations.

| Calscience l | | | | | | , In | IC. | | | **** | | | anaaninteineteine | | | (| CH/ | AIN | OF | CU | ST(| ODY | RE | COR | D |
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| | Way, Garden Grove, CA office locations: C | | | | 5494 | | | WO | ¢/LAE | 3 USE | ONLY | | | | | Date | | | | | | | | | |
| | r courier service / san contact <u>sales@calso</u> | nple drop off i | nformatior | | | | | 1 | 13. | -07 | 7-1 | 175 | j2 | | | ⁵ age_ | | | [| 01 | : | | | | |
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| SPECIAL INSTRUCTIONS: | | | | | | | | \mathbb{Q} | ň | Z | | | | | l Terra | | | | 5 | | 218.6 | No. | | | |
| | | | | | | | | $\langle S \rangle$ | <i>X</i> | C6-C44 | | | | | D | | | | 1IS 0. | 6010/747X | | \mathcal{O} | | | |
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| | l na bar | | 1 | NO. | sen | erved | Filte | H(g) | TPH(d) | 90 0 | | / MT | (826 | enate | 5035 | s (82 | ides | (808 | Ť | letals | | N | | | |
| USE SAMPLE ID | SAMF DATE | | MATRIX | OF CONT. | Unpreserved | Preserved | Field Filtered | Хтрн(g) | ЖтР | TPH [| , HqT | втех / мтве 🗆 | VOCs (8260) | Oxygenates (8260) | Prep (5035) 🗆 En Core | SVOCs (8270) | Pesticides (8081) | PCBs (8082) | РАНS П 827 0 SIM | T22 Metals | cr(VI) | Ľ | | | |
| 1 4.S. Train | 7/24/13 | | Unter | 200000000000000000000000000000000000000 | | $\overline{\mathbf{x}}$ | | $\overline{\mathbf{A}}$ | ~ | | _ | | Ń | Ť | | | | | | | Ť | | | | |
| 196024 | 7/24/13 | DRIC | holes | 1.10 | \checkmark | X | X | X | X | | | | X | | | | | | \bigtriangledown | | | \mathbf{X} | | | - |
| 2185034 MG/M | VSD 7/24/13 | DUIE | ndes | $\frac{10}{10}$ | $\overline{\mathbf{x}}$ | X | X | X | X | | | | $\frac{x}{x}$ | | | | | | \mathbf{x} | | | X | | | - |
| 260000 | 7/24/13 | 09120 | ively | 10 | | X | 1× | X | \sim | | | | $\overline{\mathbf{x}}$ | | | | | | \mathbf{x} | | | ~ | | | |
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DISTRIBUTION: White with final report, Green and Yellow to Client. Please note that pages 1 and 2 of 2 of our T/Cs are printed on the reverse side of the Green and Yellow copies respectively.

11/01/12 Revision

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•••• Page 34 of 36 Ēx. US Airbill 1752 FedEx Retrieval Copy 853l 6509 7500 0200 Express 1 From 4a Express Package Service Packages up to 150 lbs. *To most locations Sender's FedEx 7 FedEx First Overnigh Earliest next business month delivery to select locations 1 FedEx Priority Overnight 5 FedEx Standard Overnight 6 Date Account Numbe Sender's OTHO 3 FedEx 2Day Second business day FedEx Envelope ration DÂYTA 261 SOK 20 FedEx Express Saver Name FedEx Envelope rate not available. Minimum charge: One-pound rate 4b Express Freight Service Packages over 150 lbs nat Phre 83 FedEx 3Day Freight 7 FedEx 1Day Freight* 8 FedEx 2Day Freight Address Call for Confirmation Dept/Floor/Suite/Roor * Declared value limit \$500 5 Packaging 2 1 Other 2 FedEx Pak* 3 Includes FedEx Small Pek, FedEx Large Pak, and FedEx Study Pak 3 FedEx Box 4 FedEx Tube 712 FedEx Cih 61 Envelope* Your Internal Billing Reference 2 **Special Handling** 6 Include FedEx address in Section 3. 31 HOLD Saturday at FedEx Location Available ONLY for FedEx Prin Gownight and FedEx ZDay to select locations 1 HOLD Weekday at FedEx Location Not available for FedEx First Overnight SATURDAY Delivery Available ONLY for FedEx Priority Overnight, FedEx 2Day, FedEx 1Day Freight, and FedEx 2Day Freight to select ZIP codes 50 To 3 **Recipient's** 4-17) Name Perion to Select 2: 1 Covers Does this shipment contain dangerous goods? One hox must be checked. Vo 4 Yes Shipper's Declarecton Shipper's Declarecton 6 Dry Ice Dry Ice, 9, UN 1845 No 4 Yes As per attached Shipper's Declaration Company Cargo Aircraft Only Dangerous goods (including Dry Ice) cannot be shipped in FedEx packaging. Recipient's ł. Obtain Recip. 7 Payment Bill to: Address

 Bill to:
 Enter FedEx Acct. No. or Credit Card No. below.
 Obtain Hecip.

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 Recipient
 3
 Third Party
 4
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 5
 Cesh/Check

 We control dolive Dept/Four/Suite/Room Sender Acct. No. in Section 1 will be billed. 1 Address To requisive a peckage he held at a specific HedEx location, prim FedEx address here. 2 39 FedEx Accl. No. Credit Card No. Exp. Dato 84 Total Packages Fotal Weight Total Charge 3 2 FOVE State 71P N Confit Card Au Tour liability is littlifed to \$100 unless you declare a higher value. See the FodEx Service Guide for details 1200 71 650 627.24 SARANGANAMARY 4820-262951 # 18-2 SHIP DATE: 24JUL13 ACTWGT: 43.9 LB CAD: /POS1400 DIMS: 22x13x12 IN ORIGIN ID: HNLA 467 BILL RECIPIENT UNITED STATES US **TO SAMPLE CONTROL** CALSCIENCE 7440 LINCOLN WAY **GARDEN GROVE CA 92841** 0552390 (714) 895-5494 REF: DEPT FedEx Express 13111302120126 26 JUL AA FRI 2 of 2 MPS# 7958 0141 6730 2DAY ** Mstr# 8531 6209 1700 0200 92841 Z APVA CA-US SNA

| | Page 35 of 36 DRDER #: 13-07- |
|--|---|
| CLIENT: <u>Env. Science</u> | PT FORM Cooler <u> </u> of <u>2</u> DATE: <u>07 /26 / 13</u> |
| TEMPERATURE: Thermometer ID: SC3 (Criteria: 0.0 °C - 6.0 Temperature 2 3 °C - 0.2 °C (CF) = 2 6 Image: Sample(s) outside temperature criteria (PM/APM contacted by a sample(s) outside temperature criteria but received on ice/chill Image: Sample(s) outside temperature criteria but received on ice/chill Image: Sample(s) outside temperature criteria but received on ice/chill Image: Sample(s) outside temperature criteria but received on ice for transition | o°C, not frozen except sediment/tissue) oC |
| Ambient Temperature: | Initial: |
| | Not Present □ N/A Initial: Not Present Initial: |
| SAMPLE CONDITION: Chain-Of-Custody (COC) document(s) received with samples. COC document(s) received complete | n sample labels. |
| □ No analysis requested. □ Not relinquished. □ No date/time re Sampler's name indicated on COC Sample container label(s) consistent with COC Sample container(s) intact and good condition Proper containers and sufficient volume for analyses requested | |
| Analyses received within holding time pH/Res. Cl/Diss. Sulfide/Diss. Oxygen received within 15-min Proper preservation noted on COC or sample container | □ □ □ □ □ □ □ □ □ □ □ □ □ □ □ □ □ |
| Volatile analysis container(s) free of headspace Tedlar bag(s) free of condensation CONTAINER TYPE: | |
| Solid: 4ozCGJ BozCGJ 16ozCGJ Sleeve (Water: VOA VOAh VOAna2 125AGB 125AGB 500AGB 500AGJ 500AGJs 250AGB 250CGB 250PB 250PBny 125PB 125PBznna 100PJ 100 Air: Tedlar [®] Canister Other: Trip Blank Lo | □125AGBp |
| Container: C: Clear A: Amber P: Plastic G: Glass J: Jar B: Bottle Z: Ziploc/Reseal Preservative: h: HCL n: HNO ₃ na ₂ :Na ₂ S ₂ O ₃ na: NaOH p: H ₃ PO ₄ s: H ₂ SO ₄ u: Ultra-pure | lable Bag E: Envelope Reviewed by: M |

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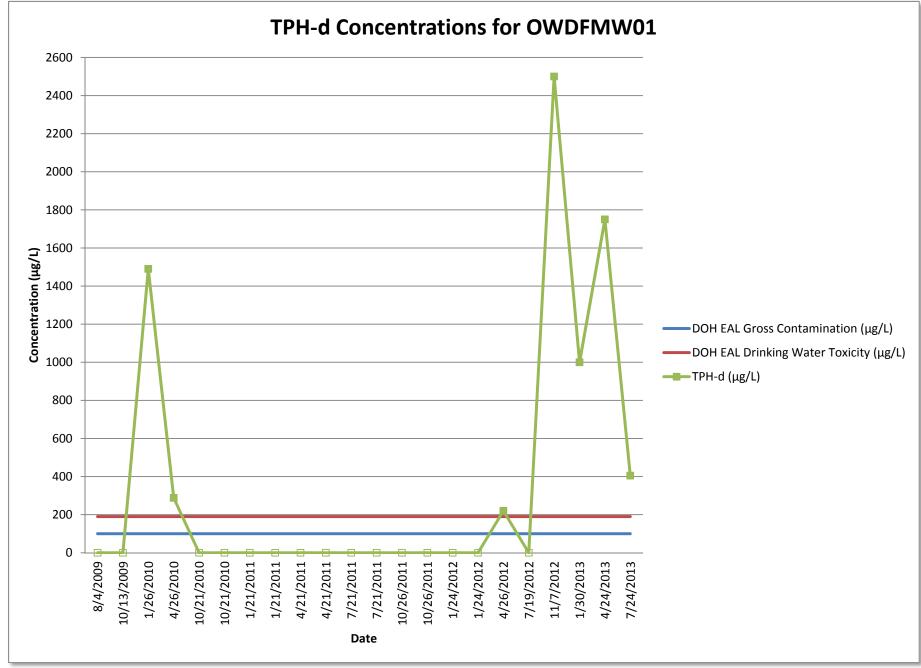
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| Laiscience Lonvironmental Laboratories, inc. | WORK ORDER | R #: 13-0 7 | 7-00 | 52 |
| SAM | PLE RECEIPT F | ORM d | cooler 2 | of 2 |
| CLIENT: Env. Science | | | 07 /26 | |
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| Temperature <u>3</u> .2°C-0.2 | °C (CF) = <u>3</u> . ∂ °C | 🛛 Blank | ☐ Sample | 4 |
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| □ Received at ambient temperature, I | | • • | | |
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| CUSTODY SEALS INTACT: | in an | | | |
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| SAMPLE CONDITION: | | Yes | No | N/A |
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| COC document(s) received complete | | AND . | | |
| Collection date/time, matrix, and/or # of co | | | | |
| □ No analysis requested. □ Not relinquis | | | | |
| Sampler's name indicated on COC Sample container label(s) consistent with | | | | |
| Sample container(s) intact and good cor | | | | |
| Proper containers and sufficient volume | | | | |
| Analyses received within holding time | | | | |
| pH/Res. Cl/Diss. Sulfide/Diss. Oxygen re | | * | | e l |
| Proper preservation noted on COC or sa | | | | |
| Unpreserved vials received for Volatiles | | | | |
| Volatile analysis container(s) free of hea | | | | Þ |
| Tedlar bag(s) free of condensation | | | | |
| Solid: □4ozCGJ □8ozCGJ □16ozC | GJ □Sleeve () □EnO | Cores [®] □Terra | Cores [®] □ | |
| Water: □VOA □VOAh □VOAna₂ □1 | 25AGB 0125AGBh 0125AG | GBp ZIAGB | □1AGB na ₂ □ | 1AGB s |
| □500AGB | 250AGB | GB s □1PB | □1PBna □ | 500PB |
| □250PB , 250PBny □125PB □125PE | Bznna □100PJ □100PJna₂ [| | □ | |
| Air: DTedlar [®] DCanister Other: D Container: C: Clear A: Amber P: Plastic G: Glass J: J Preservative: h: HCL n: HNO ₃ na ₂ :Na ₂ S ₂ O ₃ na: NaOH | Jar B: Bottle Z: Ziploc/Resealable Bag | E: Envelope | Reviewed by: | tose |

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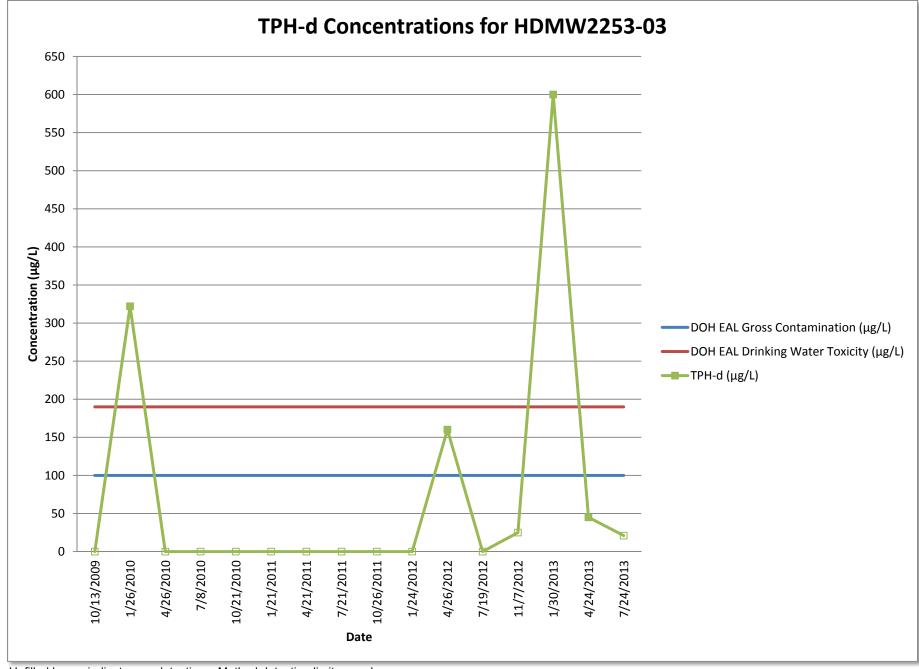
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APPENDIX D

Historical Groundwater Exceedance Trends



Data points for 10/21/2010 through 1/24/2012 and 11/07/2012 through 7/24/2013 are the average of the primary and duplicate samples. Unfilled boxes indicate non-detections. Method detection limits are shown.



Unfilled boxes indicate non-detections. Method detection limits are shown.

APPENDIX E

Waste Disposal Manifest

| | | NON-HAZARDOUS WASTE MANIFEST | 1. Generator ID Number HTR 000 | | 2. Page 1 of 1 | 3. Emergency Respor | | 4. Waste Tr | acking Nu | |
|-------------|---------------------------------------|--|--|--|----------------------------------|---|--------------------------------------|---|-------------------------|---|
| | ╞┝ | 5. Generator's Name and Mailir | na Address | | L | Generator's Site Aridin | | | | 000019559 |
| | | COMNAVREG HAW | ÀII, C/O NAVFAC ROAD, ATTN: ESTI | HAWAII, COD RELITA HIGA |)E PRJ4 | 2 RED | , | ILK FUEI | , | HIC8553-04 RAGE FACILITY |
| | | 6. Transporter 1 Company Nam | | · | | | | U.S. EPA ID | Number | |
| | | | RCIAL SERVICES, | LLC. | 80 | 08-545-4599 | ə | | | 0 0 0 9 7 8 2 4 |
| | | 7. Transporter 2 Company Nam | le | | | | | U.S. EPA ID I | Number | |
| | | 8. Designated Facility Name an | | . — | 8(| 08-682-8284 | ł | U.S. EPA ID I | <u>D 9 8</u> Number | 32443715 |
| | | 91-125 KAOMI | T SERVICES, INC. LOOP 96707 | | | | | ΗI | D 9 | 82443715 |
| | ſ | 9. Waste Shipping Name | | ······································ | | 10. Co | ntainers | 11. Total | 12. Unit | |
| | | | | | | No. | Туре | Quantity | Wt./Vol. | |
| | | (WELL PURG | L NOT REGULATED E AND DECONTAMIN | |) | 0.01 | | 00020 | G | NON-RCRA |
| ENI | 5 | 2. | | | | | | | | |
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| | | 4 | 00 PPM | PH= | 6 | | | | | |
| | ŀ | 13. Special Handling Instruction | | (c) | | | | | L | <u>ka internet i sente de la servera de la s</u> La servera de la servera de |
| | | 9b1:NR | 121 | too DDM. |) | 2008 | 9b1: | | TOTAL | HALOGEN: |
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| | | BRAKE CERAY BREOM _ HA | DOES NOT CONTAIN POBS GREA | IER HAZARDOUG MATERS | ALS AND/OR | HAZARDOUS WASTER | · | | | |
| | | 44 OFUEDATODIO CETEDOD | 'S CERTIFICATION- L bareby declar | e that the contents of this o | consignment a rding to applic | re fully and accurately d able international and n | escribed above b ational governme | by the proper shi intal regulations. | pping nam | e, and are classified, packaged, |
| | | marked and labeled/placard | ed, and are in all respects in proper (| condition for transport acco | | | | | | |
| | | T4. GENERATOR'S/OFFEROR marked and labeled/placard Generator's/Offeror's Printed/Ty Estrelito | ed, and are in all respects in proper of ped Name | condition for transport acco | Sig | Etre | lita | Hiz | $\overline{\mathbf{x}}$ | Month Day Year 08 23 3 |
| | , | marked and labeled/placard Generator's/Offeror's Printed/Ty | ed, and are in all respects in proper of ped Name | condition for transport acco | Sig Export from U | Extre | lita_ | Hiz | <u>\</u> | |
| | | marked and labeled/placard Generator's/Offeror's Printed/Ty Estrelito 15. International Shipments Transporter Signature (for expo | ed, and are in all respects in proper of ped Name Higg Import to U.S. rts only): | condition for transport acco | | J.S. Port of | lita_ entry/exit: aving U.S.: | Hiz | <u> </u> | |
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