## Quarterly Groundwater Monitoring Report Red Hill Fuel Storage Facility

Pearl Harbor, Oahu, Hawaii

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## **Executive Summary**

This Quarterly Groundwater Monitoring Report presents the results of groundwater sampling conducted on October 22, November 11, and December 16, 2008 at the United States (US) Navy Bulk Fuel Storage Facility at Red Hill, Oahu, Hawaii (the Facility). The sampling and reporting was conducted by TEC Inc. (TEC) for the Air Force Center for Engineering and the Environment (AFCEE) and the Fleet Industrial Supply Center (FISC) at Pearl Harbor, Hawaii. This report is part of a series of quarterly groundwater monitoring reports provided by the US Navy to the State of Hawaii Department of Health (HDOH) in accordance with HDOH's release response requirements. Currently, there are 18 active and 2 inactive, 12.5 million gallon, field-constructed underground storage tanks (USTs) located at the Facility.

### Background

In 2002, the United States (US) Navy installed a groundwater monitoring well (currently named RHMW01) into the basal aquifer, directly down-gradient from the Facility, within the lower access tunnel. Groundwater samples from this well indicated that petroleum from the Facility had migrated to the basal aquifer (AMEC, 2002). In 2005, the US Navy began quarterly monitoring of the aquifer to protect their down-gradient drinking water resource associated with the US Navy Well 2254-01. The US Navy Well 2254-01 is located approximately 3,000 feet downgradient from the Red Hill Fuel Storage Facility and provides approximately 24 % of the potable water to the Pearl Harbor Water System (PHWS).

By September 2005, the US Navy had installed two more groundwater monitoring wells (RHMW02 and RHMW03) within the Facility lower access tunnel, a background groundwater monitoring well (RHMW04) up-gradient from the Facility at ground surface adjacent to the US Navy Firing Range, and a groundwater monitoring well within the US Navy Well 2254-01 infiltration gallery (RHMW2254-01).

All five wells were sampled twice as part of a comprehensive environmental investigation and risk assessment (TEC, 2006). For this investigation, groundwater samples were analyzed for petroleum constituents and compared against HDOH Drinking Water Environmental Action Levels (EALs) (HDOH, 2005a). In addition, a three-dimensional (3-D) groundwater model was developed to produce site-specific risk-based action levels (SSRBLs) for compounds of concern. The results of this modeling effort indicated that Jet Propulsion (JP)–5 fuel presented the biggest risk to the US Navy water supply, due to its mobility and toxicity. Finally, the model indicated that a non-aqueous plume (free product) of JP-5 must migrate to within 1,100 feet of the US Navy Well 2254-01 infiltration gallery for HDOH EALs to be exceeded within the gallery. Based on this, free-product must be observed at RHMW01 for EALs to be exceeded at the US Navy Well 2254-01.

During the Summer and Fall of 2008, HDOH updated their EALs, which resulted in significant changes to the action levels associated with methylnaphthalenes. The drinking water toxicity EAL for these compounds were each previously 240  $\mu$ g/L, assuming they were not carcinogens. Evidence that they are carcinogenic to humans has now been accepted by the United States

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Environmental Protection Agency (USEPA), and HDOH adopted more rigorous EALs of 4.7 μg/L for 1-methylnaphthalene and 24 μg/L for 2-methylnaphthalene (HDOH, 2008).

The drinking water EAL for naphthalene was also updated during this process. Previously, HDOH based their naphthalene EAL on USEPA Region 9 Preliminary Remediation Goal (USEPA PRG) of 6.2  $\mu$ g/L, which is associated with a non-cancer Hazard Index of 1. HDOH has updated their naphthalene drinking water EAL to 17  $\mu$ g/L, in deference to the California Department of Public Health's Drinking Water Notification Levels (HDOH, 2008).

Finally, the drinking water EAL for TPH-DRO was increased from 100  $\mu$ g/L to 210  $\mu$ g/L, although the gross contamination EAL for TPH-DRO remains 100  $\mu$ g/L.

### Groundwater Protection Plan

In 2008, the US Navy completed the *Red Hill Bulk Fuel Storage Facility Final Groundwater Protection Plan* (TEC, 2008), or Plan, which specified SSRBLs for each well, and actions that would occur for the pertinent cases, based on categories for each groundwater monitoring well (Categories 1 through 4). The main object of the Plan is to protect groundwater quality entering the US Navy Well 2254-01, which provides potable water to the PHWS. This is accomplished by comparing petroleum concentrations in the Facility wells (RHMW01, RHMW02, and RHMW03) to the SSRBLs and taking the corresponding action. A secondary but important objective of the Plan is to identify leaking USTs by evaluating increasing concentration trends, or the sudden and lasting presence of free product in one or more groundwater monitoring wells. In the current quarterly groundwater monitoring report, the current water quality is compared to these categories and actions are recommended, based on those specified in the Plan.

### **Current Results**

In October 2008, four normal samples were collected from RHMW01, RHMW02, RHMW03 and the US Navy Well 2254-01, along with the required quality control samples (duplicate, matrix spike, spike duplicate, trip blank). Samples were analyzed for Total Petroleum Hydrocarbons (TPH) quantified as Diesel-Range Organics (DRO) and Gasoline Range Organics (GRO), Volatile Organic Compounds (VOCs), Polynuclear Aromatic Hydrocarbons (PAHs), and lead. Naphthalene and 1-methylnaphthalene were detected at trace concentrations in the sample from US Navy Well 2254-01; therefore, the well was re-sampled in December 2008 for VOCs. VOCs were not detected in the re-sample aliquot.

TPH-DRO was detected at 459 micrograms per liter ( $\mu$ g/L) in RHMW01, at 5,420  $\mu$ g/L (average of normal and duplicate sample) in RHMW02, and 244  $\mu$ g/L in RHMW03. The HDOH Drinking Water EAL and Site-Specific Risk Based Level for TPH-DRO are 210  $\mu$ g/L and 4,500  $\mu$ g/L, respectively. Two PAHs were detected by USEPA method 8270C SIM in the normal and duplicate samples associated with RHMW02 at average concentrations above the HDOH Drinking Water EALs: naphthalene at 89.85  $\mu$ g/L (HDOH EAL is 17  $\mu$ g/L) and 1-methylnaphthalene at 67.25  $\mu$ g/L (HDOH toxicity EAL is 4.7  $\mu$ g/L). Naphthalene was also measured by USEPA method 8260B in RHMW02, at an average concentration of 242  $\mu$ g/L.

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### Trend Analysis

At RHMW01, concentrations of TPH-DRO have been greater than the HDOH Drinking Water EAL since September 2005 but less than 25 percent of the SSRBL of 4,500  $\mu$ g/L. In October 2008, TPH-DRO increased in concentration following a decreasing trend over three previous rounds, since January 2008.

At RHMW02, concentrations of TPH-DRO have been greater than the HDOH Drinking Water EAL since September 2005 and greater than 50 percent of the SSRBL of 4,500  $\mu$ g/L over that same period. TPH-DRO has been increasing in concentration over the last four rounds, since January 2008, and the average concentration from the October 2008 sampling event was greater than the SSRBL (solubility limit). Three PAHs (naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene) also have exceeded the EALs since September 2005. In October 2008, these PAHs decreased in concentrations following increasing trends over three previous rounds (since January 2008). There are no SSRBLs set for these PAHs to date.

At RHMW03, concentrations of TPH-DRO have fluctuated around the HDOH EAL since September 2005 and are significantly lower than at RHMW01 and RHMW02. TPH-DRO at RHMW03 has been increasing in concentration over the past three rounds, since April 2008.

### Current Groundwater Status

Based on the monitoring event that occurred in October 2008, no free product was observed at RHMW01, RHMW02, or RHMW03. Results from groundwater samples indicate that RHMW01 and RHMW03 are presently in Category 2 status, since TPH-DRO is greater than the HDOH EAL for drinking water (210  $\mu$ g/L), but less than half the SSRBL of 4,500  $\mu$ g/L (solubility limit of JP-5). Category 2 response at RHMW01 and RHMW03 requires:

- 1. Quarterly reports to be sent to HDOH; and
- 2. Initiation of a leak determination program to identify if tanks are leaking.

Results from groundwater samples indicate that RHMW02 is presently in Category 4 status, since TPH-DRO (4,540  $\mu$ g/L and 6,300  $\mu$ g/L [duplicate]) is greater than the HDOH EAL for drinking water (210  $\mu$ g/L), and greater than the SSRBL of 4,500  $\mu$ g/L (solubility limit of JP-5). Category 4 response at RHMW02 requires:

- A. Send quarterly reports to HDOH
- C. Notify HDOH verbally within 1 day and follow with written notification in 30 days
- D. Notify FISC Chain of Command within 1 day
- E. Send Type 1 Report to HDOH
- F. Send Type 2 Report to HDOH
- I. Remove sampling pumps, measure product in pertinent wells with interface probe, reinstall pumps if product is not detected.
- J. Immediately evaluate tanks for leaks
- K. Collect samples from nearby Halawa Deep Monitoring Well (2253-03) and OWDF MW01

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- M. Prepare for alternative water source at US Navy Well 2254-01
- N. Re-measure for product every month with reports to HDOH

### **Conclusions and Recommendations**

There is no indication of an imminent threat to the US Navy Well 2254-01 water resources based on this report, since petroleum concentrations at RHMW01 remain less than half the SSRBLs. However, the groundwater status at RHMW02 is Category 4 since the associated TPH-DRO result was greater than the SSRBL (4,500  $\mu$ g/L). The Groundwater Protection Plan requires specific responses to Category 4, which should be conducted.

The US Navy plans to install an additional monitoring well in the Facility lower access tunnel and dedicated oil/water interface probes in four lower access tunnel wells in the coming year to better monitor contaminant migration in the basal drinking water aquifer.

It is also recommended that quarterly groundwater sampling for TPH-DRO, TPH-GRO, VOCs, PAHs, and lead be continued as previously scheduled until new data indicates a different schedule is required.

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## 1.0 Introduction

This report presents the results of the 13<sup>th</sup> groundwater sampling and analysis event, conducted in October 2008 (with follow-up sampling in November and December) at the Red Hill Fuel Storage Facility, Oahu, Hawaii (hereafter referred to as "the Facility"). The Facility consists of 18 active and 2 inactive underground storage tanks (USTs) operated by the Fleet Industrial and Supply Center (FISC), Pearl Harbor. The groundwater sampling and analysis event is part of a groundwater monitoring program for the UST site in response to past UST releases, previous environmental investigations, and recommendations from the State of Hawaii Department of Health (HDOH).

## 1.1 Project Objective

This groundwater sampling and analysis project was performed to evaluate the presence of chemicals of potential concern in groundwater underlying the Facility. The project was conducted to ensure the Navy remains in compliance with HDOH UST release response requirements as described in Hawaii Administrative Rules (HAR) 11-281 Subchapter 7, Release Response Action. The groundwater sampling and analysis procedures generally followed the procedures described in *Red Hill Bulk Fuel Storage Facility Groundwater Protection Plan* (TEC Inc [TEC], 2008) (Also referred to as "the Plan").

This groundwater sampling and analysis event was conducted by TEC under Air Force Center for Engineering and the Environment (AFCEE) Contract Number F41624-03-D-8618, Task Order 021.

## 1.2 Previous Reports

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

- 1. Groundwater Sampling Report, First Quarter 2005 (submitted April 2005);
- 2. Groundwater Sampling Report, Second Quarter 2005 (submitted August 2005);
- 3. Groundwater Sampling Report, Third Quarter 2005 (submitted November 2005);
- 4. Groundwater Sampling Report, Fourth Quarter 2005 (submitted February 2006);
- 5. Groundwater Monitoring Results, July 2006 (submitted September 2006);
- 6. Groundwater Monitoring Results, December 2006 (submitted January 2007);
- 7. Groundwater Monitoring Results, March 2007 (submitted May 2007);
- 8. Groundwater Monitoring Results, June 2007 (submitted August 2007);
- 9. Groundwater Monitoring Results, September 2007 (submitted October 2007);
- 10. Groundwater Monitoring Results, January 2008 (submitted March 2008);
- 11. Groundwater Monitoring Results, April 2008 (submitted May 2008); and
- 12. Groundwater Monitoring Results, July 2008 (submitted October 2008).

### 1.3 Background

The following sections provide a description of the site and information on the Facility and USTs.

### 1.3.1 Site Description

The Facility is located in Halawa Heights on Oahu, Hawaii. Land adjacent to the north of the Facility is occupied by Halawa Correctional Facility and private businesses. Land to the south and west of the Facility includes the Coast Guard Reservation. Moanalua Valley is located east of the Facility (Dawson, 2006).

The Navy Public Works Department operates a potable water infiltration tunnel approximately 1,550 feet hydraulically down-gradient from the Facility (Dawson, 2006). The US Navy Well 2254-01 is located approximately 3,000 feet down-gradient (west) of the Facility and provides approximately 24 % of the potable water to the Pearl Harbor Water System, which serves approximately 52,200 military consumers (TEC, 2008).

### 1.3.2 Facility Information

The Facility consists of 18 active and 2 inactive USTs operated by Navy FISC Pearl Harbor. Each UST has a capacity of 12.5 million gallons. The Facility is located approximately 100 feet above the basal aquifer (Dawson, 2006).

### 1.3.3 UST Information

The USTs were constructed in the early 1940s. The tanks were constructed of steel and currently contain Jet Propulsion (JP)–5 fuel and F-76 (diesel marine fuel). Previously, several tanks stored Navy Special Fuel Oil, Navy Distillate, aviation gasoline, and motor gasoline. Each tank measures approximately 245 feet in height and 100 feet in diameter. The upper domes of the tanks lie at depths varying between approximately 100 feet and 200 feet below the existing ground surface (TEC, 2006).

## 1.4 Previous Environmental Investigations

**1998 to 2001:** From 1998 to 2001, the Navy conducted an investigation at the Facility to assess potential releases from the fuel storage USTs and piping systems. In February 2001, the Navy installed a one-inch diameter sentinel well RHMW01 (previously known as MW-V1D) to monitor for contamination of the basal aquifer underlying the Facility. The well was installed and completed at approximately 100 feet below grade within the underground access tunnel. At the time of well completion, depth to water in RHMW01 was measured at 86 feet below grade (Dawson, 2006).

In February 2001, groundwater samples collected from sentinel well RHMW01 contained total petroleum hydrocarbons (TPH) concentrations ranging from 883 micrograms per liter ( $\mu$ g/L) to 1,050  $\mu$ g/L and total lead ranging from 10.4  $\mu$ g/L to 15  $\mu$ g/L. The maximum total lead concentration in the samples was equal to the primary drinking water standard of 15  $\mu$ g/L for lead and exceeded the HDOH Tier 1 groundwater action level of 5.6  $\mu$ g/L (Dawson, 2006).

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**2005** – **Groundwater Sampling:** The Navy began quarterly groundwater sampling at existing monitoring wells in 2005. Dawson Group, Inc. collected groundwater samples from RHMW01 and the Red Hill Navy Pump Station (US Navy Well 2254-01) in February, June, September, and December 2005.

Samples collected in February and June 2005 were not filtered in the field prior to analysis for lead. Analytical results for samples collected from RHMW01 indicated concentrations of total lead were above the HDOH Tier 1 action level of  $5.6~\mu g/L$ . The results were not considered appropriate for risk assessment since the sample had not been filtered. In addition, lead was not a component of fuels from the tanks near RHMW01. Lead may have been part of the Facility construction material (TEC, 2007).

Samples were filtered in September and December 2005, and dissolved lead concentrations were below the HDOH Tier 1 action level. Concentrations of all other contaminants of potential concern were below HDOH Tier 1 action levels.

**2005** – **Site Investigation:** As part of a site investigation, TEC installed three groundwater monitoring wells at the Facility between June and September 2005. Well RHMW02 was installed in the lower access tunnel near Tanks 5 and 6. Well RHMW03 was installed in the lower access tunnel near Tanks 13 and 14. Well RHMW04 was installed hydraulically upgradient of the USTs to provide geochemistry for water moving through the basal aquifer beneath the Facility. Wells RHMW02 and RHMW03 were completed to depths of approximately 125 feet below the tunnel floor, and well RHMW04 was completed to a depth of approximately 300 feet below ground surface outside the tunnel. Groundwater samples were collected from the three newly installed wells and two existing wells (RHMW01 and US Navy Well 2254-01) in September 2005.

Naphthalene and trichloroethylene were detected in samples collected from RHMW02 at concentrations greater than the HDOH Tier 1 action levels. Lead was detected in the sample collected from RHMW01 at a concentration greater than the HDOH Tier 1 action level; however, the sample was not filtered in the field prior to analysis. Analytical results for filtered samples obtained by Dawson during the same period indicated concentrations of dissolved lead were below the HDOH Tier 1 action level.

**2006** – **Site Investigation:** Dedicated sampling pumps were installed in five wells (RHMW01, RHMW02, RHMW03, RHMW04, and US Navy Well 2254-01). TEC collected groundwater samples from the wells in July 2006. The groundwater samples were analyzed for petroleum constituents. Naphthalene was detected in samples collected from RHMW02 at concentrations above the HDOH Tier 1 action level.

In September 2005, with concurrence from the HDOH, the Navy decided to use the newer HDOH Environmental Action Levels (EALs) for the Red Hill Site Investigation and Risk Assessment project. The EALs are current and provide action levels for more chemicals, and are much more useful for conducting screening risk assessments. Since the HDOH June 2005 Policy Letter stated that the two sets of action levels should not be mixed, the Tier 1 screening levels

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presented in HAR Section 11-281-78 would no longer be used to evaluate environmental impact at the Facility.

**2006** – **Groundwater Sampling:** Groundwater samples were collected in December 2006. Analytical results indicated the following:

- No chemicals were detected in groundwater from US Navy Well 2254-01 or RHMW03;
- TPH as diesel range organics (TPH-DRO) was detected in groundwater above the HDOH Drinking Water EALs in RHMW01; and
- TPH as gasoline range organics (TPH-GRO), TPH-DRO, and naphthalene were detected in groundwater above the HDOH Drinking Water EALs in RHMW02.

**2007** – **Groundwater Sampling:** Groundwater samples were collected in March, June, and September 2007. Analytical results indicated the following:

- No chemicals were detected above HDOH Drinking Water EALs at US Navy Well 2254-01;
- TPH-DRO exceeded HDOH Drinking Water EALs at RHMW01 during all three sampling events;
- TPH-GRO exceeded HDOH Drinking Water EALs at RHMW02 in March;
- TPH-DRO and naphthalene exceeded HDOH Drinking Water EALs at RHMW02 during all three sampling events;
- 1-methylnaphthalene and 2-methylnaphthalene exceeded HDOH Drinking Water EAL for taste and odor at RHMW02 during all three sampling events; and
- TPH-DRO exceeded HDOH Drinking Water EALs at RHMW03 in June.

**2008** – **Groundwater Sampling:** Groundwater samples were collected in January, April, and July 2008. Analytical results indicated the following:

- No chemicals were detected above HDOH Drinking Water EALs at US Navy Well 2254-01;
- TPH-DRO exceeded HDOH Drinking Water EALs at RHMW01;
- TPH-GRO did not exceed HDOH Drinking Water EALs at RHMW02;
- TPH-DRO, naphthalene, 1-methylnaphthalene, and 2-methylnaphthalene exceeded HDOH Drinking Water EALs at RHMW02; and
- TPH-DRO exceeded HDOH Drinking Water EALs at RHMW03.

## 1.5 Regulatory Updates

During the summer and fall of 2008, HDOH updated their EALs, which resulted in significant changes to the action levels associated with methylnaphthalenes. The drinking water toxicity EAL for these compounds was previously 240  $\mu$ g/L, assuming they were not carcinogens. Evidence that they are carcinogenic to humans has now been accepted by the United States Environmental Protection Agency (USEPA), and HDOH adopted more rigorous EALs of 4.7  $\mu$ g/L for 1-methylnaphthalene and 24  $\mu$ g/L for 2-methylnaphthalene, corresponding to a residential tap water scenario, and a 1 in a million cancer risk (HDOH, 2008).

The drinking water EAL for naphthalene has also been updated during this process. Previously, HDOH based their naphthalene EAL on USEPA Region 9 Preliminary Remediation Goal (USEPA PRG) of 6.2  $\mu$ g/L, which is associated with a non-cancer Hazard Index of 1. HDOH has updated their naphthalene drinking water EAL to 17  $\mu$ g/L, in deference to the California Department of Public Health's Drinking Water Notification Levels, a Hazard Index of 2.7 (HDOH, 2008).

Finally, the drinking water EAL for TPH-DRO was increased from 100  $\mu$ g/L to 210  $\mu$ g/L, although the gross contamination EAL for TPH-DRO remains 100  $\mu$ g/L.

## 2.0 Sample Collection and Analyses

The majority of field activities relating to groundwater sample collection were conducted on October 22, 2008. Groundwater samples were collected from three monitoring wells located inside the Facility lower access tunnel and one monitoring well located at the Red Hill Navy Pump Station. Sampling and analysis were conducted according to *Red Hill Bulk Fuel Storage Facility Groundwater Protection Plan* (TEC, 2008). A total of seven samples were collected; one normal sample from monitoring wells US Navy Well 2254-01, RHMW01, RHMW02 and RHMW03, one duplicate sample from RHMW02 (Sampled as RHMWA01 and reported as RHMW02-WG11D), and one matrix spike and matrix spike duplicate from US Navy Well 2254-01.

The groundwater at RHMW01 was resampled for dissolved lead on November 3, 2008 because the original sample was not filtered in the field.

Due to trace concentrations of naphthalene detected in the October 22, 2008 sample from the US Navy Well 2254-01, an additional sample was collected for Volatile Organic Compounds (VOCs) at US Navy Well 2254-01 on December 16, 2008, including one duplicate for VOCs.

## 2.1 Monitoring Well Purging

The groundwater monitoring wells were purged and sampled using a dedicated pump system. Well purging was considered complete when no less than three successive water quality parameter measurements had stabilized within approximately 10 percent. Field parameters were measured at regular intervals during well purging and included pH, temperature, specific conductivity, dissolved oxygen, and turbidity. Purge water was disposed in the Facility oil/water separator system.

## 2.2 Groundwater Sample Collection

Each monitoring well was sampled immediately following purging. All wells were sampled directly from their dedicated bladder pump system. Samples were placed into sampling containers with appropriate preservatives (i.e., hydrochloric acid [HCl] for volatile organic analysis, nitric acid [HNO<sub>3</sub>] for dissolved lead). Dissolved lead samples were filtered in the field and placed in preserved bottles. Sample containers were labeled with the date, sample identification number, type of analysis, and sampler's name. The containers were placed on ice in sample coolers and transported under chain-of-custody procedures to the certified laboratory for analysis.

## 2.3 Groundwater Sample Analyses

Groundwater samples were analyzed by SGS Environmental Service, Inc. in Anchorage, Alaska for TPH-DRO and TPH-GRO by EPA Method 8015B, VOCs by EPA Method 8260B, polynuclear aromatic hydrocarbons (PAHs) by EPA Method 8270C SIM, and dissolved lead by EPA Method 6020.

## 3.0 Groundwater Sample Analytical Results

This section provides a summary of analytical results for groundwater samples collected from three monitoring wells located in the lower access tunnel of the Facility and one monitoring well located at the Red Hill Navy Pump Station. Duplicate sample results from monitoring wells RHMW02 and RHMW2254-01 are reported in this document as RHMW02D and RHMW2254-01D, respectively. A summary of groundwater analytical results is included in Table 1. Complete analytical laboratory reports are provided in Appendix A.

## 3.1 October 2008 Sample Analytical Results

All groundwater samples were analyzed for TPH-DRO, TPH-GRO, VOCs, PAHs, and dissolved lead. Naphthalene and 1-methylnaphthalene were detected at US Navy Well 2254-01 at 0.05F  $\mu g/L$  and 0.03F  $\mu g/L$ , respectively. This is below the HDOH Drinking Water EALs of 17  $\mu g/L$  for naphthalene and 4.7  $\mu g/L$  for 1-methylnaphthlene. Data qualifier "F" indicates the result is between the method detection limit (MDL) and the reporting limit (RL) and considered an estimated value. However, naphthalene or 1-methynapthalene are generally not observed in groundwater from JP-5 releases when TPH is not detected. For this reason, these low-concentration PAH results may not be from the groundwater, but the source of the expected cross-contamination is not known. Following the receipt of the preliminary results, FISC implemented an immediate re-sampling of the US Navy Well 2254-01 for VOCs, including naphthalene.

### RHMW01

Four petroleum constituents were detected at RHMW01: TPH-DRO, fluorene, naphthalene, and lead. TPH-DRO was detected at RHMW01 at 459  $\mu$ g/L, which is above the HDOH Drinking Water EAL of 210  $\mu$ g/L. All other petroleum constituents were below HDOH Drinking Water EALs at RHMW01.

### RHMW02

Eight petroleum constituents were detected at RHMW02: TPH-DRO, TPH-GRO, 1-methylnaphthalene, 2-methylnaphthalene, acenaphthene, fluorene, naphthalene, benzene, and ethylbenzene. TPH-DRO was detected at RHMW02 in the normal and duplicate samples, at 4,540  $\mu$ g/L and 6,300  $\mu$ g/L, respectively. Both results exceeded the HDOH EAL and site-specific risk based level (SSRBL) (the HDOH EAL is 210  $\mu$ g/L and the SSRBL is 4,500  $\mu$ g/L).

Naphthalene was analyzed by USEPA Method 8270C SIM and USEPA Method 8260B. USEPA Method 8260B produced the highest naphthalene concentrations, which averaged 242  $\mu$ g/L from the normal and duplicate sample (HDOH EAL is 17  $\mu$ g/L). In addition, 1-methylnaphthalene and 2-methylnaphthalene were detected by USEPA Method 8270C SIM in the normal and duplicate samples. The average result for 1-methylnaphthalene was 67.25  $\mu$ g/L, greater than the HDOH toxicity EAL of 4.7  $\mu$ g/L. The average result for 2-methylnaphthalene was 13.2  $\mu$ g/L, less than the HDOH toxicity EAL of 24  $\mu$ g/L but greater than the HDOH taste and odor EAL of 10  $\mu$ g/L. All other petroleum constituents were below HDOH Drinking Water EALs at RHMW02.

RHMW03

## Four constituents were detected at RHMW03: TPH-DRO, 1-methylnaphthylene, 2-methylnaphthalene, and naphthalene. TPH-DRO was detected at RHMW03 at 244 µg/L, slightly above the HDOH EAL of 210 µg/L. All other petroleum constituents were below HDOH

Drinking Water EALs at RHMW03.

### US Navy Well 2254-01 as RHMW2254

RHMW2254 was sampled twice during this reporting period. On October 22, 2008, trace concentrations (less than the reporting limit but greater than the detection limit) of naphthalene (0.05  $\mu$ g/L) and 1-methylnaphthalene (0.03  $\mu$ g/L) were detected by USEPA Method 8270C SIM. On December 16, 2008, RHMW2254 was re-sampled by USEPA Method 8260B. Although no petroleum constituents were observed in the re-sample, the method detection limit for naphthalene was 0.62  $\mu$ g/L and 1-methylnaphthalene was not analyzed.

Table 1. Analytical Results for Quarterly Groundwater Monitoring Release Response Report (October 22, 2008 and December 16, 2008) Red Hill Fuel Storage Facility, Pearl Harbor, Hawaii

PAHs- VOCs- VOCy- UG/L- Q- Q- U- F- 200-													8260B (VOCs)																		8270C SIM (PAHs)						8015B (Petroleum)	Method
PAHs - Polynuclear aromatic hydrocarbons OCs - Volatile organic compounds IG/L - Micrograms per Liter Q - Data qualifier U - Indicates that the compound was analyzed for but not detected at or above the stated limit F - Indicates that the compound was identified but the concentration was above the MDL and below the RL F - Indicates that the compound was identified but the concentration was above the MDL and below the RL Toxicity-based environmental action levels, Table D-1b, Screening For Environmental Concerns At Sites With Contaminated Soil and Groundwater, HDOH, 2008 Taste, odor and solubility thresholds, Table G-1, Screening For Environmental Concerns At Sites With Contaminated Soil and Groundwater, HDOH, 2008	LEAD	VINYL CHLORIDE	trans-1,2-DICHLOROETHENE TRICHLOROETHYLENE (TCE)	TOLUENE	STYRENE TETRACHLOROETHYLENE(PCE)	NAPHTHALENE	METHYL ISOBUTYL KETONE	METHYL ETHYL KETONE (2-BUTANONE)	M BLAY ENE (SIM OF ISOMEDS)	DIBROMOMETHANE	cis-1,2-DICHLORGETHYLENE	CHLOROFORM	CHLOROETHANE	CARBON FEIRACHLORIDE CHLOROBENZENE	BROMOMETHANE	BROMOFORM	BROMODICHLOROMETHANE	ACETONE	1,3-DICHLOROBENZENE	1,3-DICHLOROBENZENE	1,2-DICHLOROPROPANE	1,2,4-TRICHLOROBENZENE	1,1-DICHLOROETHANE	1,1,2,2-TETRACHLOROETHANE 1.1.2-TRICHLOROETHANE	1,1,1,2-TETRACHLOROETHANE 1,1,1-TRICHLOROETHANE	PHENANTHRENE PYRENE	NAPHTHALENE	FLUORENE INDENO(1,2,3-c,d)PYRENE	FLUORANTHENE	DIRENZ/2 HANTHR ACENE	BENZO(k)FLUORANTHENE	BENZO(b)FLUORANTHENE	BENZO(a)ANTHRACENE BENZO(a)PYRENE	ANTHRACENE	ACENAPHTHENE	1-METHYLNAPHTHALENE	TPH as GASOLINE RANGE ORGANICS	Chemical
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### 3.2 Groundwater Contaminant Trend

Groundwater samples have been collected and analyzed by TEC since September 2005. Figure 1 shows TPH trends in groundwater at the Facility. Figure 2 shows PAH trends in groundwater at the Facility. In these figures, open icons (without fill) represent the method detection limit for chemicals that were not detected.

### RHMW01

TPH-GRO has only been detected in April 2008 at 13.6F  $\mu$ g/L, which is below the HDOH Drinking Water EAL of 100  $\mu$ g/L. In October 2008, TPH-GRO was not detected. TPH-DRO was detected above the HDOH Drinking Water EAL during all groundwater sampling events and showed a decreasing trend until the January 2008 sampling event. Concentration of TPH-DRO observed during the October 2008 sampling event was greater than the concentrations observed in April and July 2008, but lower than the January 2008 sampling event.

### RHMW02

TPH-GRO was detected in nine of ten sampling events since September 2005, and exceeded the HDOH Drinking Water EAL three times during 2006 and 2007. The maximum concentration detected was 148  $\mu$ g/L. The concentration of TPH-DRO was relatively stable at RHMW02 until January 2008, ranging from 2,250 to 3,180  $\mu$ g/L. However, since January 2008 it has shown an increasing trend with concentrations up to 5,420  $\mu$ g/L, well above the HDOH EAL, and above the SSRBL of 4,500  $\mu$ g/L. PAHs at RHMW02 remain above the HDOH Drinking Water EALs, and concentrations have decreased in October 2008 after showing an increasing trend since January 2008.

### RHMW03

TPH-GRO has never been detected. TPH-DRO had shown a decreasing trend until the January 2008 sampling event. Concentrations of TPH-DRO observed during the October 2008 sampling event were higher than the previous rounds. In general, concentrations of petroleum-related compounds at RHMW03 have been the lowest of the three wells located within the Facility.

## 3.3 Results of Oil/Water Interface Measurements

The presence and thickness of light-non aqueous phased liquids (LNAPL), otherwise known as free product, released from the USTs were measured during this sampling event using a 300 ft Heron Oil/Water Interface Meter. The static water levels were measured to a precision of  $\pm$  0.01 ft and fuel thickness was measured to a precision of  $\pm$  0.01 ft with this equipment.

In January 2008, fuel was measured in monitoring wells RHMW01 and RHMW02 at a thickness of < 0.01 ft, but has not been observed in other monitoring wells as of this document.

Measurements to determine the presence and thickness of fuel were conducted at RHMW01, RHMW02, and RHMW03 during the current sampling round. No fuel product was observed in any of these wells during this event (see Table 2).

**Table 2. Oil/Water Interface Measurements** 

	RHM	IW01	RHM	IW02	RHMW03			
Date	SWL (ft)	LNAPL (ft)	SWL (ft)	LNAPL (ft)	SWL (ft)	LNAPL (ft)		
January 2008	17.74	< 0.01	18.78	< 0.01	NT	NT		
July 2008	19.04	0.00	18.91	0.00	18.86	0.00		
October 2008	18.61	0.00	18.56	0.00	18.82	0.00		
November 2008	18.50	0.00	18.45	0.00	18.51	0.00		

**SWL** Static water level, elevation above mean sea level

**LNAPL** Light Non-Aqueous Phased Liquid, fuel product on groundwater

ft FeetNT Not Taken

## 3.4 Groundwater Status

Compounds of concern are defined as those petroleum-related chemicals that have been observed in the groundwater samples above the HDOH EALs. In accordance with the *Red Hill Bulk Fuel Storage Facility Final Groundwater Protection Plan* (TEC, 2008), Table 3 defines the constituents of concern in groundwater at the Facility and the SSRBLs and updated EALs for each (HDOH 2008).

**Table 3. Action Levels for Constituents of Concern** 

Chemical	EAL (µg/L)	SSRBL (µg/L)
<b>Petroleum Mixtures</b>		
TPH-DRO	210	4,500
TPH-GRO	100	4,500
<b>Semi-Volatile Compounds</b>		
1-Methylnaphthalene	4.7	NA
2-Methylnaphthalene	24	NA
Naphthalene	17	NA

NA – Not applicable or not determined

SSRBLs are applicable at RHMW01, RHMW02, and RHMW03

EALs are applicable at US Navy Well 2254-01

In addition, the Plan defines four results categories of groundwater status for the Facility, based on concentrations of constituents of concern in RHMW01, RHMW02, RHMW03 and the US Navy Well 2254-01, and requires specific responses when these categories are observed during quarterly groundwater sampling. Table 4 describes each of the four results categories and identifies response actions to be taken in accordance with the Plan.

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**Table 4. Results Categories and Response Actions to Changes in Groundwater Status** 

Results Category	RHMW02 or RHMW03	RHMW01	US Navy Pumping Well 2254-01
Results Category 1: Result above detection limit but below drinking water EAL and trend for all compounds stable or decreasing	A	A	A,D,M,E,P
Results Category 2: Trend for any compound increasing or drinking water EAL exceeded	A, B	A, B	A,B,C,D,E,F,G,K, L,O
Results Category 3: Result Between 1/10X SSRBL and SSRBL for benzene, or between 1/2X SSRBL and SSRBL for TPH	A,B,G,H,I,J	A,B,E,G,H,I,J	A,B,C,D,E,F,G,I,J, K,L,O
Results Category 4: Result Exceeding any SSRBL or petroleum product observed	A,C,D,E,F,I,J, K,M,N	A,C,D,E,F,I, J,K,M,N,O	A,C,D,E,F,G,I,J,K, L,O

### Specific Responses:

- A. Send quarterly reports to HDOH
- B. Begin program to determine the source of leak
- C. Notify HDOH verbally within 1 day and follow with written notification in 30 days
- D. Notify FISC Chain of Command within 1 day
- E. Send Type 1 Report (see box below) to HDOH
- F. Send Type 2 Report (see box below) to HDOH
- G. Increase monitoring frequency to once per month (if concentrations increasing)
- H. Notify HDOH verbally within 7 days and follow with written notification in 30 days
- I. Remove sampling pumps, measure product in pertinent wells with interface probe, re-install pumps if product is not detected.
- J. Immediately determine leaking tank
- K. Collect samples from nearby Halawa Deep Monitoring Well (2253-03) and OWDF MW01
- L. Provide alternative water source at 2254-01
- M. Prepare for alternative water source at US Navy Well 2254-01
- N. Re-measure for product every month with reports to HDOH
- O. Install additional monitoring well downgradient

### **Report Types**

**HDOH Type 1 Report** 

- Re-evaluate Tier 3 Risk Assessment/groundwater model results
- Proposal to HDOH on a course of action

### **HDOH Type 2 Report**

• Proposal for groundwater treatment

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Based on the monitoring event that occurred in October and November 2008, no free product was observed at RHMW01, RHMW02, or RHMW03. Results from groundwater samples indicate that RHMW01 and RHMW03 are presently in Category 2 status, since TPH-DRO is greater than the HDOH EAL for drinking water (210  $\mu$ g/L), but less than half the SSRBL of 4,500  $\mu$ g/L (solubility limit of JP-5). Category 2 response at RHMW01 and RHMW03 requires:

- 1. Quarterly reports to be sent to HDOH; and
- 2. Initiation of a leak determination program to identify if tanks are leaking.

Results from groundwater samples indicate that RHMW02 is presently in Category 4 status, since TPH-DRO (4,540  $\mu$ g/L and 6,300  $\mu$ g/L [duplicate]) is greater than the SSRBL of 4,500  $\mu$ g/L (solubility limit of JP-5). Category 4 response at RHMW02 requires:

- A. Send quarterly reports to HDOH
- C. Notify HDOH verbally within 1 day and follow with written notification in 30 days
- D. Notify FISC Chain of Command within 1 day
- E. Send Type 1 Report to HDOH
- F. Send Type 2 Report to HDOH
- I. Remove sampling pumps measure product in pertinent wells with interface probe, re-install pumps if product is not detected.
- J. Immediately evaluate tanks for leaks
- K. Collect samples from nearby Halawa Deep Monitoring Well (2253-03) and OWDF MW01. For permission to sample 2253-03, call DLNR Commission on Water Resource Management (808) 587-0214, DLNR.CWRM@Hawaii.gov
- M. Prepare for alternative water source at US Navy Well 2254-01
- N. Re-measure for product every month with reports to HDOH

In response to the previous Category 3 conditions at RHMW02, oil/water interface measurements were collected in October 2008 and November 2008 at RHMW01, RHMW02 and RHMW03. To date, there is no evidence of fuel on groundwater at any of these wells based on oil/water interface measurements.

Category 4 groundwater status at RHMW02 requires additional efforts, including:

- Re-evaluate risk assessment and groundwater model (TEC, 2007) to ensure both are valid and protective of human health and the environment under the existing conditions and provide a report describing the results of this re-evaluation to the HDOH;
- Evaluate potential requirement for groundwater treatment and provide a report describing the results of this re-evaluation to the HDOH;
- Implement monthly oil/water interface measurements at RHMW01, RHMW02, RHMW03, and provide monthly letter reports of the results;
- Evaluate tanks associated with the middle section of the Facility (Tanks 5, 6, 7, 8, 9, 10, 11, 12, 13, and 14);
- Collect samples from nearby Halawa Deep Monitoring Well (2253-03) and OWDF MW01;
- Prepare for alternative water source at US Navy Well 2254-01, as necessary.

## 4.0 Summary and Conclusions

At RHMW02, the concentration of TPH-DRO in groundwater exceeded the SSRBL of 4,500 μg/L, which indicates Category 4 groundwater status at RHMW02.

- The SSRBL of TPH-DRO is set at the solubility limit of JP-5 and exceeding the SSRBL implies that JP-5 is in direct contact with groundwater in the vicinity of RHMW02.
- Oil/water interface measurements were collected in October 2008 and November 2008 from RHMW01, RHMW02, and RHMW03 and no fuel thickness was measured.
- The concentration of TPH-DRO measured at RHMW01 in October 2008 was one tenth
  of the SSRBL. RHMW01 is down-gradient from RHMW02 and between RHMW02 and
  the US Navy Well 2254-01, an important drinking water source for the Pearl Harbor
  Water System.
- The US Navy Well 2254-01 is not imminently threatened at this time; however, conditions should be monitored closely to determine if any USTs in the Facility are currently leaking fuel into the subsurface.
- Category 4 activities should be implemented, including:
  - Re-evaluate risk assessment and groundwater model (TEC, 2007) to ensure both are valid and protective of human health and the environment under the existing conditions and provide a report describing the results of this re-evaluation to the HDOH;
  - Evaluate potential requirement for groundwater treatment and provide a report describing the results of this re-evaluation to the HDOH;
  - Implement monthly oil/water interface measurements at RHMW01, RHMW02, RHMW03, and provide monthly letter reports of the results;
  - Evaluate tanks associated with the middle section of the Facility (Tanks 5, 6, 7, 8, 9, 10, 11, 12, 13, and 14);
  - Collect samples from nearby Halawa Deep Monitoring Well (2253-03) and OWDF MW01;
  - Prepare for alternative water source at US Navy Well 2254-01, as necessary.

The US Navy plans to install an additional monitoring well (RHMW05) in the lower access tunnel of the Facility between RHMW01 and the US Navy Well 2254-01 to better monitor the quality of the groundwater moving from the Facility to the US Navy Well 2254-01.

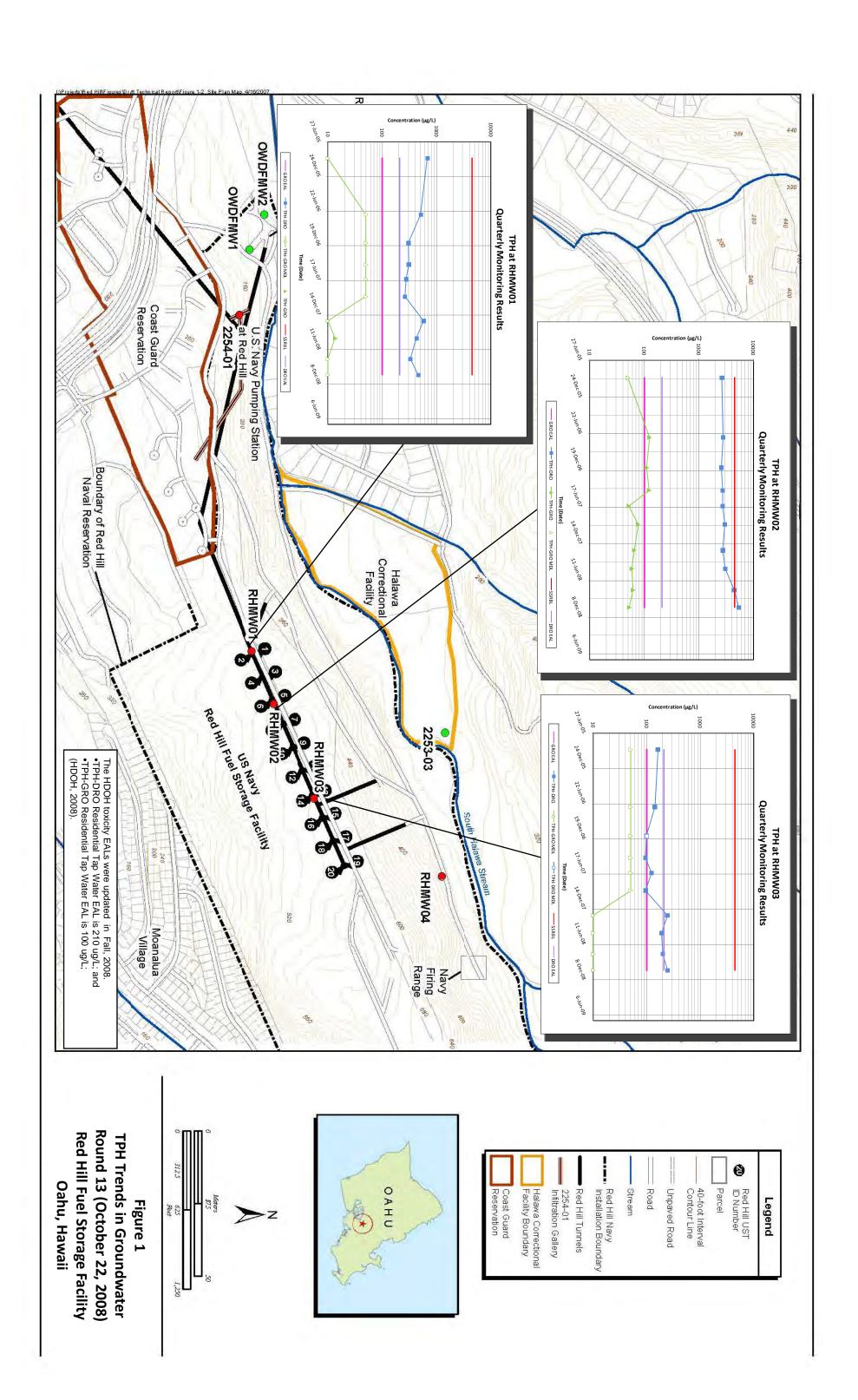
Although the results from the groundwater sample collected from US Navy Well 2254-01 in October 2008 showed trace concentrations of naphthalene and 1-methylnaphthalene, no other petroleum constituents were detected in this sample. This well was re-sampled in December 2008. The sample was analyzed by USEPA Method 8620B and no petroleum constituents were detected.

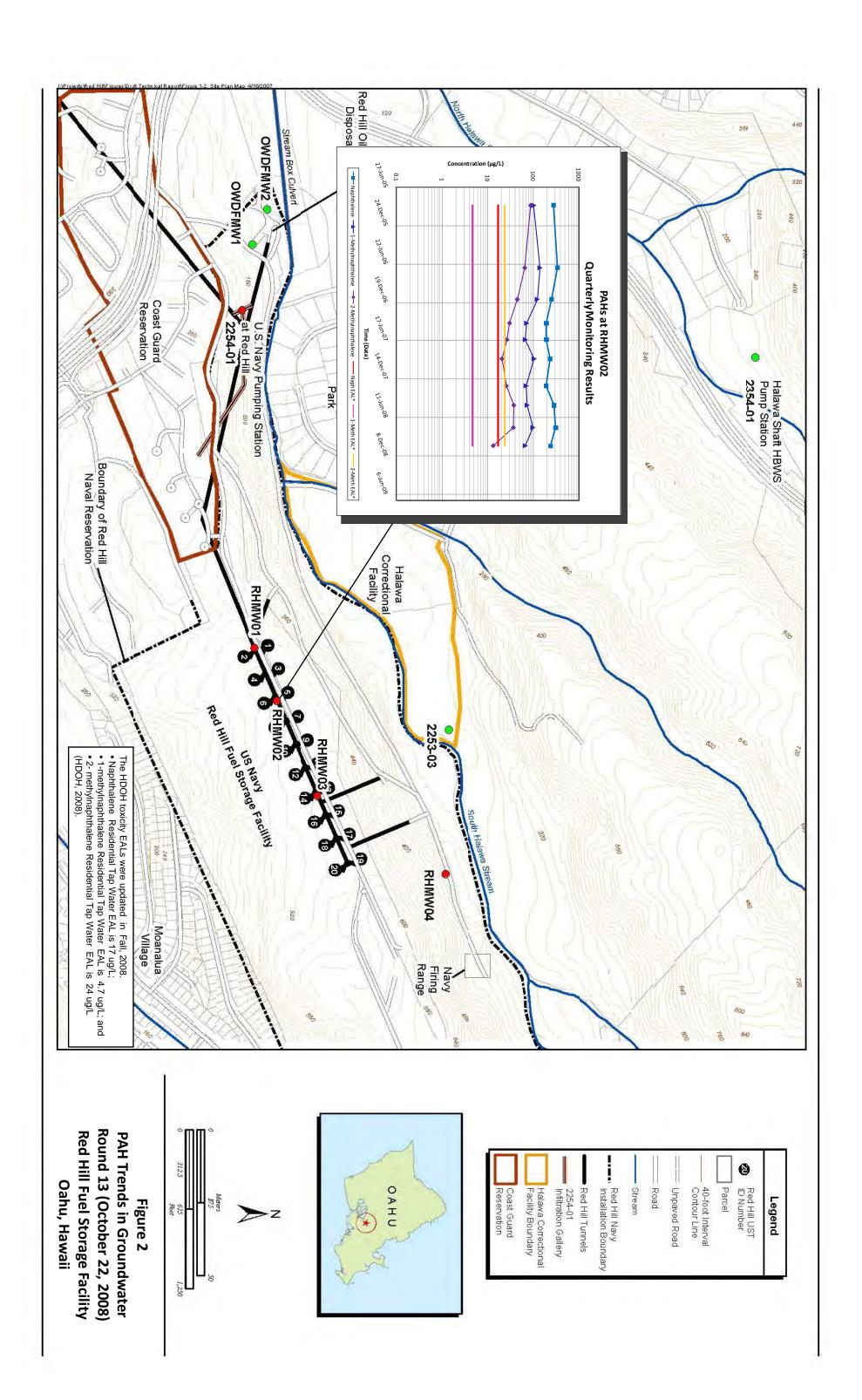
It is recommended that RHMW01, RHMW02, RHMW03, and RHMW05 be evaluated monthly for the presence of fuel on groundwater, in accordance with the Plan. The US Navy plans to install dedicated oil/water interface probes in each of these wells to facilitate these measurements.

results.

It is also recommended that quarterly groundwater sampling for TPH-DRO, TPH-GRO, VOCs, PAHs, and lead be continued as previously scheduled until new data indicates that a different schedule is required. The quarterly collection and analysis of groundwater samples will continue to monitor the quality of the groundwater located beneath the Facility. Groundwater monitoring

reports will be submitted to the HDOH upon receipt and evaluation of laboratory analytical





## 5.0 References

AMEC. 2002. *Red Hill Bulk Fuel Storage Facility Investigation Report*, Prepared for NAVFAC Pacific, August 2002.

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The Environmental Company, Inc. and AMEC. 2005. Red Hill Bulk Fuel Storage Facility Work Plan, Pearl Harbor, Hawaii. June.

TEC Inc. 2006. Red Hill Bulk Fuel Storage Facility, Final – Addendum Planning Documents, Pearl Harbor, Hawaii. May.

TEC Inc. 2007. Red Hill Bulk Fuel Storage Facility, Final Technical Report, Pearl Harbor, Hawaii. August.

TEC Inc. 2008. Red Hill Bulk Fuel Storage Facility, Final Groundwater Protection Plan, Pearl Harbor, Hawaii. January.

# Appendix A Laboratory Analytical Reports



# SGS Environmental Services Alaska Division Level II Laboratory Data Report

Project: 9121-003 Red Hill BFSF

Client: The Environmental Company, Inc. (TEC)

SGS Work Order: 1085813

Released by:

### Contents:

Cover Page
Case Narrative
Final Report Pages
Quality Control Summary Forms
Chain of Custody/Sample Receipt Forms

### Note:

Unless otherwise noted, all quality assurance/quality control criteria is in compliance with the standards set forth by the proper regulatory authority, the SGS Quality Assurance Program Plan, and the National Environmental Accreditation Conference.

### SGS Environmental Services Inc.

### **Case Narrative**

Customer: THEENVC The Environmental Company, Inc. (TEC)

Project: 1085813 9121-003 Red Hill BFSF

NPDL WO:

Refer to the sample receipt form for information on sample condition.

### 1085813001 PS RHMW2254-WG13

8015C DRO - LCS recovery is outside of acceptance criteria (biased low). The sample was re-extracted outside of the 14 day holding time and reanalyzed. The results did confirm with passing QC. Results for all associated samples are considered estimated.

### 1085813004 PS RHMW03-WG13

8015C DRO - MS/MSD were not spiked due to lab error. See LCS/LCSD for precision and accuracy.

### 1085813005 PS RHMW02-WG13

8015C DRO - The pattern is consistent with a weathered middle distillate.

8015C DRO - MS/MSD were not spiked due to lab error. See LCS/LCSD for precision and accuracy.

### 1085813006 PS RHMWA01-WG13

8015C DRO - The pattern is consistent with a weathered middle distillate.

8015C DRO - MS/MSD were not spiked due to lab error. See LCS/LCSD for precision and accuracy.

### 1085813007 PS RHMW01-WG13

8015C DRO - Unknown hydrocarbon with several peaks is present.

8015C DRO - MS/MSD were not spiked due to lab error. See LCS/LCSD for precision and accuracy.

### 1085813002 BMS RHMW2254-WG13 MS

8015C DRO - LCS recovery is outside of acceptance criteria (biased low). The sample was re-extracted outside of the 14 day holding time and reanalyzed. The results did confirm with passing QC. Results for all associated samples are considered estimated.

### 868545 MS 1085813001MS

8015C DRO - MS/MSD were not spiked due to lab error. See LCS/LCSD for precision and accuracy.

### 1085813003 BMSD RHMW2254-WG13 MSD

8015C DRO - LCS recovery is outside of acceptance criteria (biased low). The sample was re-extracted outside of the 14 day holding time and reanalyzed. The results did confirm with passing QC. Results for all associated samples are considered estimated.

### 868546 MSD 1085813001MSD

8015C DRO - MS/MSD were not spiked due to lab error. See LCS/LCSD for precision and accuracy.

#### 869198 LCS XXX/203131

8015C DRO - LCS recovery is outside of acceptance criteria (biased low). Results for all associated samples are considered estimated.



## Laboratory Analytical Report

Client: The Environmental Company, Inc.

1001 Bishop Street Ste 1400

**ASB Tower** 

Honolulu, HI 96813

Attn: Jeff Hart

T: (808)528-1445 F:(808)528-0768

jshart@tecinc.com

Project: 9121-003 Red Hill BFSF

Workorder No.: 1085813

### Certification:

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, other than the conditions noted on the sample data sheet(s) and/or the case narrative. This certification applies only to the tested parameters and the specific sample(s) received at the laboratory.

If you have any questions regarding this report, or if we can be of further assistance, please contact your SGS Project Manager.

Tamara Rentz tamara.rentz@sgs.com Project Manager



Print Date: 11/20/2008

Enclosed are the analytical results associated with this workorder.

As required by the state of Alaska and the USEPA, a formal Quality Assurance/Quality Control Program is maintained by SGS. A copy of our Quality Assurance Plan (QAP), which outlines this program is available at your request.

The Laboratory certification numbers are AK971-05 (DW), UTS-005 (CS) and AK00971 (Micro) for ADEC and AK100001 for NELAP (RCRA methods: 1020A, 1311, 6010B, 7470A, 7471A, 9040B, 9045C, 9056, 9060, 8015B, 8021B, 8081A/8082, 8260B, 8270C).

Except as specifically noted, all statements and data in this report are in conformance to the provisions set forth by the SGS QAP, the National Environmental Laboratory Accreditation Program and, when applicable, other regulatory authorities.

If you have any questions regarding this report or if we can be of any assistance, please contact your SGS Project Manager at 907-562-2343.

The following descriptors may be found on your report which will serve to further qualify the data.

MDL Method Detection Limit

PQL Practical Quantitation Limit (reporting limit).

CL Control Limit

U Indicates the analyte was analyzed for but not detected. F Indicates value that is greater than or equal to the MDL.

J The quantitation is an estimation.

ND Indicates the analyte is not detected

B Indicates the analyte is found in a blank associated with the sample.

\* The analyte has exceeded allowable regulatory or control limits.

D The analyte concentration is the result of dilution.

GT Greater Than LT Less Than

Q QC parameter out of acceptance range.

M A matrix effect was present.

E The analyte result is above the calibrated range.

R Rejected

DF Analytical Dilution Factor

JL The analyte was positively identified, but the quantitation is a low estimation.

<Surr> Surrogate QC spiked standard

<Surr/IS> Surrogate / Internal Standard QC spiked standard

QC Quality Control
QA Quality Assurance
MB Method Blank

LCS (D) Laboratory Control Sample (Duplicate)

MS(D) Matrix Spike (Duplicate)
BMS(D) Site Specific Matrix Spike
RPD Relative Percent Difference
ICV Initial Calibration Verification
CCV Continuous Calibration Verification
MSA Method of Standard Addition

Notes: Soil samples are reported on a dry weight basis unless otherwise specified

All DRO/RRO analysese are integrated per SOP.



### SAMPLE SUMMARY

Print Date: 11/20/2008

Client Name: The Environmental Company, Inc. (TEC)

Project Name: 9121-003 Red Hill BFSF

Workorder No.: 1085813

### **Analytical Methods**

Lab Sample ID

**Analytical Method** Method Description 8270 PAH SIM Semi-Vol GC/MS Liq/Liq ext. 8270D SIMS AFCEE 3.1 8260 (W) SW8260B Dissolved Metals by ICP-MS SW6020 DRO by 8015B (W) SW8015C GRO (W) SW8015C

Client Sample ID

### Sample ID Cross Reference

1085813001	RHMW2254-WG13
1085813002	RHMW2254-WG13 MS
1085813003	RHMW2254-WG13 MSD
1085813004	RHMW03-WG13
1085813005	RHMW02-WG13
1085813006	RHMWA01-WG13
1085813007	RHMW01-WG13
1085813008	TB01-WG13



Client Sample ID: RHMW2254-WG13

SGS Ref. #: 1085813001

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 09:30 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

Analyst: NRB

### Dissolved Metals by ICP/MS

<u>Parameter</u>	Result	PQL/CL	MDL_	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> Batch	Qualifiers
Lead	ND	1.00	0.310	ug/L	5	MMS5719	MXX2103	1
Batch Information								
Analytical Batch: MMS5719		Prep Batch: M	IXX21031			Initial Prep	Wt./Vol.: 50	mL
Analytical Method: SW6020		Prep Method:	SW3010A			Prep Extrac	t Vol.: 50 m	L
Analysis Date/Time: 11/14/08 20:12		Prep Date/Tim	ne: 11/04/08 18:	15		Container II	D:10858130	01-G



Client Sample ID: RHMW2254-WG13

SGS Ref. #: 1085813001

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 09:30 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Fuels Department**

<u>Parameter</u>	<u>Result</u>	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> Batch	Qualifiers
Gasoline Range Organics	ND	100	10.0	ug/L	1	VFC9248	VXX1896	7
4-Bromofluorobenzene <surr></surr>	99.1	50-150		%	1	VFC9248	VXX1896	7
Batch Information								
Analytical Batch: VFC9248		Prep Batch	: VXX18967			Initial Prep	Wt./Vol.: 5 n	nL
Analytical Method: SW8015C		Prep Metho	d: SW5030B			Prep Extra	ct Vol.: 5 mL	
Analysis Date/Time: 10/30/08 10:09		Prep Date/	Γime: 10/30/08 (	08:00		Container I	D:10858130	01-A
Dilution Factor: 1						Analyst: HI	М	



Client Sample ID: RHMW2254-WG13

SGS Ref. #: 1085813001

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 09:30 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Semivolatile Organic Fuels Department**

						<u>Analytical</u>	<u>Prep</u>	
<u>Parameter</u>	<u>Result</u>	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	<u>Batch</u>	<u>Batch</u>	<u>Qualifiers</u>
Diesel Range Organics	ND	0.421	0.0842	mg/L	1	XFC8338	XXX2031	3
5a Androstane <surr></surr>	93.1	50-150	0.00.2	//g/2 %	1	XFC8338	XXX2031	
Batch Information								
Daten information								
Analytical Batch: XFC8338		Prep Batch	: XXX20313			Initial Prep	Wt./Vol.: 95	0 mL
Analytical Method: SW8015C		Prep Metho	d: SW3520C			Prep Extra	ct Vol.: 1 ml	_
Analysis Date/Time: 11/05/08 09:36		Prep Date/	Time: 11/03/08 1	7:15		Container I	D:10858130	001-l
Dilution Factor: 1		·				Analyst: Gl	_	



Client Sample ID: RHMW2254-WG13

SGS Ref. #: 1085813001

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 09:30 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

						Analytical	<u>Prep</u>
<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Benzene	ND	0.400	0.120	ug/L	1	VMS10242	VXX18954
Toluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Ethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
n-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dichloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Isopropyltoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
n-Propylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Styrene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromomethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,4-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Acetone	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dibromo-3-chloropropane	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
Methyl-t-butyl ether	ND	5.00	1.50	ug/L	1	VMS10242	VXX18954
Tetrachloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromochloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3-Dichloropropane	ND	0.400	0.120	ug/L	1	VMS10242	VXX18954
1,2-Dibromoethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Carbon tetrachloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,1,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
Chloroform	ND	1.00	0.300	ug/L	1	VMS10242	VXX18954
Bromobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,3-Trichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromomethane	ND	3.00	0.940	ug/L	1	VMS10242	VXX18954
Bromochloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Vinyl chloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dichlorodifluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954

SGS Environmental Services Inc. Alaska Division 200 West Potter Drive Anchorage Alaska 99518

† (907) 562 2343 † (907) 561 5301 www.us.sgs.com



Client Sample ID: RHMW2254-WG13

SGS Ref. #: 1085813001

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 09:30 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

Barrantar	Paguit	DOI /CI	MDI	11.26.		<u>Analytical</u>	Prep
Parameter Obligation at the control of the control	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Chloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
sec-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromodichloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,1-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
2-Butanone (MEK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
Methylene chloride	ND	5.00	1.00	ug/L	1	VMS10242	VXX18954
Trichlorofluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
P & M -Xylene	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
Naphthalene	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
o-Xylene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromoform	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1-Chlorohexane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
tert-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1-Dichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
2-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Trichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
2,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Hexachlorobutadiene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Isopropylbenzene (Cumene)	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichloroethane-D4 <surr></surr>	105	73-120		%	1	VMS10242	VXX18954
Toluene-d8 <surr></surr>	104	80-120		%	1	VMS10242	VXX18954
4-Bromofluorobenzene <surr></surr>	111	76-120		%	1	VMS10242	VXX18954
					-		

### **Batch Information**

Prep Batch: VXX18954 Initial Prep Wt./Vol.: 5 mL Analytical Batch: VMS10242 Analytical Method: SW8260B Prep Method: SW5030B Prep Extract Vol.: 5 mL Analysis Date/Time: 10/27/08 17:17 Prep Date/Time: 10/27/08 11:41 Container ID:1085813001-D

Dilution Factor: 1 Analyst: JDB



Client Sample ID: RHMW2254-WG13

SGS Ref. #: 1085813001

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 09:30 Receipt Date/Time: 10/24/08 10:50

### Polynuclear Aromatics GC/MS

<u>Parameter</u>	<u>Result</u>	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Analytical Batch	Prep Batch Qualifiers
Acenaphthylene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Acenaphthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Fluorene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Phenanthrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Anthracene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Fluoranthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Pyrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo(a)Anthracene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Chrysene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[b]Fluoranthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[k]fluoranthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[a]pyrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Indeno[1,2,3-c,d] pyrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Dibenzo[a,h]anthracene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[g,h,i]perylene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Naphthalene	0.0466 J	0.100	0.0310	ug/L	1	XMS4761	XXX20276
1-Methylnaphthalene	0.0276 J	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
2-Methylnaphthalene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Terphenyl-d14 <surr></surr>	92.2	50-135		%	1	XMS4761	XXX20276

### **Batch Information**

Analytical Batch: XMS4761 Analytical Method: 8270D SIMS Analysis Date/Time: 10/30/08 14:48

Dilution Factor: 1

Prep Batch: XXX20276 Prep Method: SW3520C Prep Date/Time: 10/28/08 10:10 Initial Prep Wt./Vol.: 1000 mL Prep Extract Vol.: 1 mL Container ID:1085813001-J Analyst: JDH

Print Date: 11/20/2008



Client Sample ID: RHMW03-WG13

SGS Ref. #: 1085813004

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 11:15 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### Dissolved Metals by ICP/MS

<u>Parameter</u>	<u>Result</u>	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> Batch	Qualifiers
Lead	ND	1.00	0.310	ug/L	5	MMS5719	MXX2103	1
Batch Information								
Analytical Batch: MMS5719		Prep Batch: N	MXX21031			Initial Prep	Nt./Vol.: 50 ı	mL
Analytical Method: SW6020		Prep Method:	SW3010A			Prep Extrac	t Vol.: 50 ml	L
Analysis Date/Time: 11/14/08 22:05		Prep Date/Tir	ne: 11/04/08 18	:15		Container II	D:108581300	04-G
Dilution Factor: 5						Analyst: NF	RB	



Client Sample ID: RHMW03-WG13

SGS Ref. #: 1085813004

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 11:15 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Fuels Department**

<u>Parameter</u>	<u>Result</u>	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> Batch	Qualifiers
Gasoline Range Organics	ND	100	10.0	ug/L	1	VFC9248	VXX1896	7
4-Bromofluorobenzene <surr></surr>	94.8	50-150		%	1	VFC9248	VXX1896	7
Batch Information								
Analytical Batch: VFC9248		Prep Batch	: VXX18967			Initial Prep	Wt./Vol.: 5 n	nL
Analytical Method: SW8015C		Prep Metho	od: SW5030B			Prep Extra	ct Vol.: 5 mL	
Analysis Date/Time: 10/30/08 11:18		Prep Date/	Time: 10/30/08	08:00		Container I	D:10858130	004-A
Dilution Factor: 1						Analyst: HI	M	



Client Sample ID: RHMW03-WG13

SGS Ref. #: 1085813004

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 11:15 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

Analytical

Prep

### **Semivolatile Organic Fuels Department**

<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Batch	Batch	<u>Qualifiers</u>
Diesel Range Organics	0.244 J	0.442	0.0884	mg/L	1	XFC8323	XXX2029	0
5a Androstane <surr></surr>	101	50-150		%	1	XFC8323	XXX2029	0
Batch Information								
Analytical Batch: XFC8323		Prep Batch	n: XXX20290			Initial Prep	Wt./Vol.: 90	5 mL
Analytical Method: SW8015C		Prep Metho	od: SW3520C			Prep Extra	ct Vol.: 1 mL	
Analysis Date/Time: 10/31/08 12:33		Prep Date/	Time: 10/29/08 1	8:00		Container I	D:10858130	04-H
Dilution Factor: 1						Analyst: G	L	



Client Sample ID: RHMW03-WG13

SGS Ref. #: 1085813004

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 11:15 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

	-					Analytical	Prep
<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Benzene	ND	0.400	0.120	ug/L	1	VMS10242	VXX18954
Toluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Ethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
n-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dichloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Isopropyltoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
n-Propylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Styrene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromomethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,4-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Acetone	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dibromo-3-chloropropane	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
Methyl-t-butyl ether	ND	5.00	1.50	ug/L	1	VMS10242	VXX18954
Tetrachloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromochloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3-Dichloropropane	ND	0.400	0.120	ug/L	1	VMS10242	VXX18954
1,2-Dibromoethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Carbon tetrachloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,1,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
Chloroform	ND	1.00	0.300	ug/L	1	VMS10242	VXX18954
Bromobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,3-Trichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromomethane	ND	3.00	0.940	ug/L	1	VMS10242	VXX18954
Bromochloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Vinyl chloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dichlorodifluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954

SGS Environmental Services Inc. Alaska Division 200 West Potter Drive Anchorage Alaska 99518

† (907) 562 2343 † (907) 561 5301 www.us.sgs.com



Client Sample ID: RHMW03-WG13

SGS Ref. #: 1085813004

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 11:15 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

## **Volatile Gas Chromatography/Mass Spectroscopy**

Barrantar	Passilt	DOL/CI	***	11.26.	<b>D</b> E	<u>Analytical</u>	Prep
Parameter Obligation of the parameter	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Chloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
sec-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromodichloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,1-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
2-Butanone (MEK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
Methylene chloride	ND	5.00	1.00	ug/L	1	VMS10242	VXX18954
Trichlorofluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
P & M -Xylene	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
Naphthalene	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
o-Xylene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromoform	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1-Chlorohexane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
tert-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1-Dichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
2-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Trichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
2,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Hexachlorobutadiene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Isopropylbenzene (Cumene)	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichloroethane-D4 <surr></surr>	104	73-120		%	1	VMS10242	VXX18954
Toluene-d8 <surr></surr>	103	80-120		%	1	VMS10242	VXX18954
4-Bromofluorobenzene <surr></surr>	113	76-120		%	1	VMS10242	VXX18954
					-		

#### **Batch Information**

Prep Batch: VXX18954 Initial Prep Wt./Vol.: 5 mL Analytical Batch: VMS10242 Analytical Method: SW8260B Prep Method: SW5030B Prep Extract Vol.: 5 mL Analysis Date/Time: 10/27/08 17:50 Prep Date/Time: 10/27/08 11:41 Container ID:1085813004-D

Dilution Factor: 1 Analyst: JDB

SGS Environmental Services Inc.

Alaska Division 200 West Potter Drive Anchorage Alaska 99518 t (907) 562-2343 f (907) 561-5301 www.us.sgs.com



Client Sample ID: RHMW03-WG13

SGS Ref. #: 1085813004

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 11:15 Receipt Date/Time: 10/24/08 10:50

### Polynuclear Aromatics GC/MS

<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	DF	Batch	Batch	Qualifiers
Acenaphthylene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	1
Acenaphthene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	;
Fluorene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	1
Phenanthrene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	1
Anthracene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	;
Fluoranthene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	1
Pyrene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	i
Benzo(a)Anthracene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	;
Chrysene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	i
Benzo[b]Fluoranthene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	i
Benzo[k]fluoranthene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	i
Benzo[a]pyrene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	i
Indeno[1,2,3-c,d] pyrene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	i
Dibenzo[a,h]anthracene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	;
Benzo[g,h,i]perylene	ND	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	i
Naphthalene	0.219	0.103	0.0320	ug/L	1	XMS4761	XXX20276	i
1-Methylnaphthalene	0.0658	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	i
2-Methylnaphthalene	0.0937	0.0515	0.0155	ug/L	1	XMS4761	XXX20276	i
Terphenyl-d14 <surr></surr>	89.4	50-135		%	1	XMS4761	XXX20276	;

#### **Batch Information**

Prep Batch: XXX20276 Analytical Batch: XMS4761 Analytical Method: 8270D SIMS Prep Method: SW3520C Analysis Date/Time: 10/30/08 16:29 Prep Date/Time: 10/28/08 10:10 Dilution Factor: 1

Prep Extract Vol.: 1 mL Container ID:1085813004-J

Initial Prep Wt./Vol.: 970 mL

Print Date: 11/20/2008

Analyst: JDH

Analytical

Prep



Client Sample ID: RHMW02-WG13

SGS Ref. #: 1085813005

Dilution Factor: 5

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:45 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

Analyst: NRB

### Dissolved Metals by ICP/MS

<u>Parameter</u>	<u>Result</u>	PQL/CL	MDL_	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> Batch	<u>Qualifiers</u>
Lead	ND	1.00	0.310	ug/L	5	MMS5719	MXX2103	1
Batch Information								
Analytical Batch: MMS5719		Prep Batch: M	1XX21031			Initial Prep	Wt./Vol.: 50	mL
Analytical Method: SW6020	Prep Method: SW3010A					Prep Extract Vol.: 50 mL		
Analysis Date/Time: 11/14/08 22:12	Prep Date/Time: 11/04/08 18:15					Container II	D:10858130	05-G



Client Sample ID: RHMW02-WG13

SGS Ref. #: 1085813005

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:45 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Fuels Department**

<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> <u>Batch</u>	Qualifiers
Gasoline Range Organics	52.8 J	100	10.0	ug/L	1	VFC9248	VXX1896	7
4-Bromofluorobenzene <surr></surr>	129	50-150		%	1	VFC9248	VXX1896	7
Batch Information								
Analytical Batch: VFC9248		Prep Batch	: VXX18967			Initial Prep	Wt./Vol.: 5 r	mL
Analytical Method: SW8015C		Prep Metho	od: SW5030B			Prep Extra	ct Vol.: 5 mL	_
Analysis Date/Time: 10/30/08 11:36		Prep Date/	Time: 10/30/08	08:00		Container I	D:10858130	005-A
Dilution Factor: 1						Analyst: HI	M	



Client Sample ID: RHMW02-WG13

SGS Ref. #: 1085813005

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:45 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Semivolatile Organic Fuels Department**

						<u>Analytical</u>	Prep	
<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	<u>Batch</u>	<u>Batch</u>	<u>Qualifiers</u>
Diesel Range Organics	4.54	0.400	0.0800	mg/L	1	XFC8323	XXX2029	0
5a Androstane <surr></surr>	97.8	50-150		%	1	XFC8323	XXX2029	0
Batch Information								
Analytical Batch: XFC8323		Prep Batch	: XXX20290			Initial Prep	Wt./Vol.: 100	00 mL
Analytical Method: SW8015C		Prep Metho	od: SW3520C			Prep Extrac	ct Vol.: 1 mL	
Analysis Date/Time: 10/31/08 12:42		Prep Date/	Time: 10/29/08 1	8:00		Container I	D:10858130	05-H
Dilution Factor: 1						Analyst: Gl	_	



Client Sample ID: RHMW02-WG13

SGS Ref. #: 1085813005

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:45 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

## **Volatile Gas Chromatography/Mass Spectroscopy**

Policy   P	Columb Car Strain Graph James Special Lips						Analytical	Prep
Toluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 Ethylbenzene 0.450 J 1.00 0.310 ug/L 1 VMS10242 VXX18954 1.4-Dichlorobenzene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1.4-Dichlorobenzene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1.2-Dichloroethane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1.2-Dichloroethane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1.3.5-Trimethylbenzene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Chlorotoluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Chlorotoluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Chlorobenzene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 4-Methyl-2-pentanone (MIBK) ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Methyl-2-pentanone (MIBK) ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Isopropyltoluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Isopropyltoluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Isopropyltoluene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 4-Isopropyltoluene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 5-13-J0-Dichloropropene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 5-13-Dichloropropene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 5-13-Dichloropropene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 5-13-Dichloropropene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 5-13-Dichloro	<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>		
Toluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 Ethylbenzene 0.450 J 1.00 0.310 ug/L 1 VMS10242 VXX18954 1.4-Dichlorobenzene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1.4-Dichlorobenzene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1.2-Dichloroethane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1.2-Dichloroethane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1.3.5-Trimethylbenzene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Chlorotoluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Chlorotoluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Chlorobenzene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 4-Methyl-2-pentanone (MIBK) ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Methyl-2-pentanone (MIBK) ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Isopropyltoluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Isopropyltoluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 4-Isopropyltoluene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 4-Isopropyltoluene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 4-Isopropyltoluene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 5-13-Dichloropropene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 5-13-Dichloropropene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 5-13-Dichloropropene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 5-14-Trichlorobenzene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 5-								
Ethylbenzene	Benzene	0.140 J	0.400		ug/L	1	VMS10242	VXX18954
n-Butylbenzene         9.65         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,4-Dichlorobenzene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dichlorocherbane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,3,5-Trimethylbenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Chlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Methyl-2-pentanone (MIBK)         ND         10.0         0.310         ug/L         1         VMS10242         VXX18954           4-Methyl-2-pentanone (MIBK)         ND         10.0         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           d-Isopropyltoluene         ND         0.500         0.150	Toluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,4-Dichlorobenzene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dichloroethane         ND         0.500         0.160         ug/L         1         VMS10242         VXX18954           1,3,5-Trimethylbenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Chlorotoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Methyl-2-pentanone (MIBK)         ND         10.0         3.10         ug/L         1         VMS10242         VXX18954           4-Methyl-2-pentanone (MIBK)         ND         10.0         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropyltoluene         ND         0.0500         0.150	Ethylbenzene	0.450 J	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,3,5-Trimethylbenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Chlorotoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chlorobenzene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           4-Methyl-2-pentanone (MIBK)         ND         10.0         3.10         ug/L         1         VMS10242         VXX18954           4-Methyl-2-pentanone (MIBK)         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropytoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropytoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropytoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           ois-1,3-Dichioropropene         ND         1.00         0.310	n-Butylbenzene	9.65	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,3,5-Trimethylbenzene         ND         1,00         0.310         ug/L         1         VMS10242         VXX18954           4-Chlorotoluene         ND         1,00         0.310         ug/L         1         VMS10242         VXX18954           4-Chlorobenzene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           4-Methyl-2-pentanone (MIBK)         ND         10.0         3.10         ug/L         1         VMS10242         VXX18954           d-is-13-Dichloroethene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           d-Isopropyllouene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           d-Isopropyllouene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           d-Isopropyllouene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           d-Isopropyllouene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           n-Propyllouene         ND         0.500         0.310 <t< td=""><td>1,4-Dichlorobenzene</td><td>ND</td><td>0.500</td><td>0.150</td><td>ug/L</td><td>1</td><td>VMS10242</td><td>VXX18954</td></t<>	1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
Chlorotoluene	1,2-Dichloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
Chlorobenzene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 4-Methyl-2-pentanone (MIBK) ND 10.0 3.10 ug/L 1 VMS10242 VXX18954 cis-1,2-Dichloroethene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 d-Isopropyltoluene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 cis-1,3-Dichloropropene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 cis-1,3-Dichloropropene ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 n-Propylbenzene 12.4 1.00 0.310 ug/L 1 VMS10242 VXX18954 Styrene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 Dibromomethane ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 trans-1,3-Dichloropropene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 1,2,4-Trichlorobenzene ND 1.00 0.310 ug/L 1 VMS10242 VXX18954 1,2,2-Tetrachloroethane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1,2,2-Tetrachloroethane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1,2-Dibromo-3-chloropropane ND 2.00 0.620 ug/L 1 VMS10242 VXX18954 1,2-Dibromo-3-chloropropane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 Dibromochloromethane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 Dibromochloromethane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1,3-Dichloropropane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1,2-Dibromochlane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1,2-Dibromochlane ND 0.500 0.150 ug/L 1 VMS10242 VXX18954 1,1-Dichloropropane ND 0.500 0.310 ug	1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Methyl-2-pentanone (MIBK)         ND         10.0         3.10         ug/L         1         VMS10242         VXX18954           cis-1,2-Dichloroethene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropylloluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           6:s-1,3-Dichloropropene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           n-Propylbenzene         12.4         1.00         0.310         ug/L         1         VMS10242         VXX18954           Styrene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromomethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           12,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,4-Trichlorobenzene         ND         1.00         0.310	4-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
cis-1,2-Dichloroethene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           cis-1,3-Dichloropropene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Styrene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Styrene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromomethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,1,2,2-Tetrachloroethane         ND         0.500         0.150	Chlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
A-Isopropyltoluene	4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
cis-1,3-Dichloropropene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           n-Propylbenzene         12.4         1.00         0.310         ug/L         1         VMS10242         VXX18954           Styrene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromomethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Acetone         ND         1.00         3.10         ug/L         1         VMS10242         VXX18954           1,2-Dibromo-3-chloroptane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dibromo-3-chloroptopane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dibromo-chloromethane         ND         1.00         0.310	cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
N-Propylbenzene   12.4   1.00   0.310   ug/L   1 VMS10242 VXX18954	4-Isopropyltoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Styrene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromomethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Acetone         ND         10.0         3.10         ug/L         1         VMS10242         VXX18954           1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dibromo-3-chloropropane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Tetrachloroethane         ND         1.00         0.310         u	cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
Styrene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromomethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Acetone         ND         10.0         3.10         ug/L         1         VMS10242         VXX18954           1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dibromo-3-chloropropane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Tetrachloroethane         ND         1.00         0.310         u	n-Propylbenzene	12.4	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromomethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Acetone         ND         10.0         3.10         ug/L         1         VMS10242         VXX18954           1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Tetrachloroethene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromochloromethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310	Styrene	ND	1.00	0.310		1	VMS10242	VXX18954
1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Acetone         ND         10.0         3.10         ug/L         1         VMS10242         VXX18954           1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Tetrachloroethene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Tetrachloroethene         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Tetrachloropropane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,3-Dichloropropane         ND         0.400         0.120         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310	Dibromomethane	ND	1.00	0.310	-	1	VMS10242	VXX18954
1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Acetone         ND         10.0         3.10         ug/L         1         VMS10242         VXX18954           1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Tetrachloroethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310	trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Tetrachloroethene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromochloromethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,3-Dichloropropane         ND         0.400         0.120         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Carbon tetrachloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300	1,2,4-Trichlorobenzene	ND	1.00	0.310	-	1	VMS10242	VXX18954
1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Tetrachloroethene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromochloromethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,3-Dichloropropane         ND         0.400         0.120         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310 <td< td=""><td>Acetone</td><td>ND</td><td>10.0</td><td>3.10</td><td>-</td><td>1</td><td>VMS10242</td><td>VXX18954</td></td<>	Acetone	ND	10.0	3.10	-	1	VMS10242	VXX18954
1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1         VMS10242         VXX18954           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Tetrachloroethene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromochloromethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,3-Dichloropropane         ND         0.400         0.120         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Carbon tetrachloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Carbon tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310         <	1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1         VMS10242         VXX18954           Tetrachloroethene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Dibromochloromethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,3-Dichloropropane         ND         0.400         0.120         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Carbon tetrachloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chloromethane         ND         1.00         0.310         ug/L	1,2-Dibromo-3-chloropropane	ND	2.00	0.620	-	1	VMS10242	VXX18954
Dibromochloromethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           1,3-Dichloropropane         ND         0.400         0.120         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Carbon tetrachloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromochloromethane         ND         3.00         0.940         ug/L </td <td>Methyl-t-butyl ether</td> <td>ND</td> <td>5.00</td> <td>1.50</td> <td>-</td> <td>1</td> <td>VMS10242</td> <td>VXX18954</td>	Methyl-t-butyl ether	ND	5.00	1.50	-	1	VMS10242	VXX18954
1,3-Dichloropropane         ND         0.400         0.120         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Carbon tetrachloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromoethane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L <t< td=""><td>Tetrachloroethene</td><td>ND</td><td>1.00</td><td>0.310</td><td>ug/L</td><td>1</td><td>VMS10242</td><td>VXX18954</td></t<>	Tetrachloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,3-Dichloropropane         ND         0.400         0.120         ug/L         1         VMS10242         VXX18954           1,2-Dibromoethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Carbon tetrachloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromoethlane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         <	Dibromochloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dibromoethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Carbon tetrachloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromomethane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Bromochloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	1,3-Dichloropropane	ND	0.400	0.120	-	1	VMS10242	VXX18954
Carbon tetrachloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromomethane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	1,2-Dibromoethane	ND	1.00	0.310		1	VMS10242	VXX18954
1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1         VMS10242         VXX18954           Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromomethane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	Carbon tetrachloride	ND	1.00	0.310	-	1	VMS10242	VXX18954
Chloroform         ND         1.00         0.300         ug/L         1         VMS10242         VXX18954           Bromobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromomethane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Bromochloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	1,1,1,2-Tetrachloroethane	ND	0.500	0.150	=	1	VMS10242	VXX18954
Bromobenzene         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Chloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromomethane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Bromochloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	Chloroform	ND	1.00	0.300	-	1	VMS10242	VXX18954
Chloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromomethane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Bromochloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	Bromobenzene	ND	1.00	0.310	=	1	VMS10242	VXX18954
1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Bromomethane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Bromochloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	Chloromethane	ND	1.00	0.310	-	1		VXX18954
Bromomethane         ND         3.00         0.940         ug/L         1         VMS10242         VXX18954           Bromochloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	1,2,3-Trichloropropane	ND			-	1		VXX18954
Bromochloromethane         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954           Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	Bromomethane	ND	3.00	0.940	•	1		VXX18954
Vinyl chloride         ND         1.00         0.310         ug/L         1         VMS10242         VXX18954	Bromochloromethane	ND	1.00	0.310	-	1		VXX18954
	Vinyl chloride	ND	1.00	0.310		1		VXX18954
	Dichlorodifluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954

SGS Environmental Services Inc. Alaska Division 200 West Potter Drive Anchorage Alaska 99518

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Client Sample ID: RHMW02-WG13

SGS Ref. #: 1085813005

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:45 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

	,					Analytical	Prep	
<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	<u>Batch</u>	Batch Qualifi	ers
Chloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
sec-Butylbenzene	8.48	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Bromodichloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954	
1,1-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
2-Butanone (MEK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954	
Methylene chloride	ND	5.00	1.00	ug/L	1	VMS10242	VXX18954	
Trichlorofluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
P & M -Xylene	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954	
Naphthalene	239	40.0	12.4	ug/L	20	VMS10254	VXX18976	
o-Xylene	0.450 J	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Bromoform	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1-Chlorohexane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
tert-Butylbenzene	1.29	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1-Dichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
2-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Trichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
2,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Hexachlorobutadiene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Isopropylbenzene (Cumene)	6.64	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2-Dichloroethane-D4 <surr></surr>	104	73-120		%	1	VMS10242	VXX18954	
Toluene-d8 <surr></surr>	104	80-120		%	1	VMS10242	VXX18954	
4-Bromofluorobenzene <surr></surr>	104	76-120		%	1	VMS10242	VXX18954	



Client Sample ID: RHMW02-WG13

SGS Ref. #: 1085813005

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:45 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

						<u>Analytical</u>	<u>Prep</u>		
<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	<u>Batch</u>	<u>Batch</u>	<b>Qualifiers</b>	
Batch Information									
Analytical Batch: VMS10242		Prep Batch	: VXX18954		Initial Prep Wt./Vol.: 5 mL				
Analytical Method: SW8260B		Prep Method: SW5030B				Prep Extract Vol.: 5 mL			
Analysis Date/Time: 10/27/08 18:24		Prep Date/Time: 10/27/08 11:41				Container ID:1085813005-E			
Dilution Factor: 1						Analyst: JD	В		
Analytical Batch: VMS10254		Prep Batch	: VXX18976			Initial Prep	Nt./Vol.: 5	mL	
Analytical Method: SW8260B		Prep Method: SW5030B				Prep Extract Vol.: 5 mL			
Analysis Date/Time: 10/31/08 08:38		Prep Date/Time: 10/31/08 01:28 Container ID:1085813					D:1085813	005-C	
Dilution Factor: 20						Analyst: DS	SH		



Client Sample ID: RHMW02-WG13

SGS Ref. #: 1085813005

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:45 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### Polynuclear Aromatics GC/MS

Parameter	<u>Result</u>	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Analytical Batch	Prep Batch Qualifiers
Accomplete	ND	0.0521	0.0156		4	VN40.4704	XXX20276
Acenaphthona				ug/L	1	XMS4761	
Acenaphthene	0.365	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Fluorene Phenanthrene	0.214 ND	0.0521 0.0521	0.0156 0.0156	ug/L	1	XMS4761	XXX20276 XXX20276
	ND ND	0.0521		ug/L	1	XMS4761	
Anthracene			0.0156	ug/L	1	XMS4761	XXX20276
Fluoranthene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Pyrene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Benzo(a)Anthracene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Chrysene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Benzo[b]Fluoranthene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Benzo[k]fluoranthene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Benzo[a]pyrene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Indeno[1,2,3-c,d] pyrene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Dibenzo[a,h]anthracene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Benzo[g,h,i]perylene	ND	0.0521	0.0156	ug/L	1	XMS4761	XXX20276
Naphthalene	97.4	5.21	1.61	ug/L	50	XMS4763	XXX20276
1-Methylnaphthalene	72.1	2.60	0.781	ug/L	50	XMS4763	XXX20276
2-Methylnaphthalene	13.7	1.04	0.313	ug/L	20	XMS4763	XXX20276
Terphenyl-d14 <surr></surr>	89.6	50-135		%	1	XMS4761	XXX20276
Batch Information							
Analytical Batch: XMS4761		Prep Batch	: XXX20276			Initial Prep	Wt./Vol.: 960 mL
Analytical Method: 8270D SIMS		Prep Metho	od: SW3520C			Prep Extrac	ct Vol.: 1 mL
Analysis Date/Time: 10/30/08 17:02		Prep Date/	Time: 10/28/08 1	0:10		Container II	D:1085813005-J
Dilution Factor: 1						Analyst: JD	H
Analytical Batch: XMS4763		Prep Batch	: XXX20276			Initial Prep	Wt./Vol.: 960 mL
Analytical Method: 8270D SIMS		Prep Metho	od: SW3520C			Prep Extrac	t Vol.: 1 mL
Analysis Date/Time: 10/31/08 11:49		Prep Date/	Time: 10/28/08 1	0:10		Container II	D:1085813005-J
Dilution Factor: 20						Analyst: JD	H
Analytical Batch: XMS4763		Prep Batch	: XXX20276			Initial Prep	Wt./Vol.: 960 mL
Analytical Method: 8270D SIMS		Prep Metho	od: SW3520C			Prep Extrac	t Vol.: 1 mL
Analysis Date/Time: 10/31/08 15:57		Prep Date/	Time: 10/28/08 1	0:10		Container II	D:1085813005-J
Dilution Factor: 50						Analyst: JD	H



Client Sample ID: RHMWA01-WG13

SGS Ref. #: 1085813006

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:05 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### Dissolved Metals by ICP/MS

<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> Batch	Qualifiers
Lead	ND	1.00	0.310	ug/L	5	MMS5719	MXX2103	1
Batch Information								
Analytical Batch: MMS5719		Prep Batch:	MXX21031			Initial Prep	Wt./Vol.: 50	mL
Analytical Method: SW6020		Prep Metho	d: SW3010A			Prep Extra	t Vol.: 50 m	L
Analysis Date/Time: 11/14/08 22:19		Prep Date/1	ime: 11/04/08 1	8:15		Container I	D:10858130	06-G
Dilution Factor: 5						Analyst: NF	RB	



Client Sample ID: RHMWA01-WG13

SGS Ref. #: 1085813006

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:05 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Fuels Department**

<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> Batch	Qualifiers
Gasoline Range Organics	52.9 J	100	10.0	ug/L	1	VFC9248	VXX1896	7
4-Bromofluorobenzene <surr></surr>	130	50-150		%	1	VFC9248	VXX1896	7
Batch Information								
Analytical Batch: VFC9248		Prep Batch	: VXX18967			Initial Prep	Wt./Vol.: 5 n	nL
Analytical Method: SW8015C		Prep Metho	od: SW5030B			Prep Extra	ct Vol.: 5 mL	
Analysis Date/Time: 10/30/08 11:54		Prep Date/	Time: 10/30/08	08:00		Container I	D:10858130	006-A
Dilution Factor: 1						Analyst: HI	M	



Client Sample ID: RHMWA01-WG13

SGS Ref. #: 1085813006

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:05 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Semivolatile Organic Fuels Department**

<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Analytical Batch	Prep Batch	Qualifiers
Diesel Range Organics	6.30	0.491	0.0982	mg/L	1	XFC8323	XXX2029	00
5a Androstane <surr></surr>	104	50-150		%	1	XFC8323	XXX2029	0
Batch Information								
Analytical Batch: XFC8323		Prep Batch	: XXX20290			Initial Prep	Wt./Vol.: 81	5 mL
Analytical Method: SW8015C		Prep Metho	od: SW3520C			Prep Extra	ct Vol.: 1 mL	_
Analysis Date/Time: 10/31/08 12:52		Prep Date/	Time: 10/29/08 1	8:00		Container I	D:10858130	006-H
Dilution Factor: 1						Analyst: Gl	_	



Client Sample ID: RHMWA01-WG13

SGS Ref. #: 1085813006

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:05 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

	pp					Analytical	Prep
<u>Parameter</u>	<u>Result</u>	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Benzene	0.150 J	0.400	0.120	ug/L	1	VMS10242	VXX18954
Toluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Ethylbenzene	0.420 J	1.00	0.310	ug/L	1	VMS10242	VXX18954
n-Butylbenzene	9.47	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dichloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Isopropyltoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
n-Propylbenzene	12.5	1.00	0.310	ug/L	1	VMS10242	VXX18954
Styrene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromomethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,4-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Acetone	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dibromo-3-chloropropane	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
Methyl-t-butyl ether	ND	5.00	1.50	ug/L	1	VMS10242	VXX18954
Tetrachloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromochloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3-Dichloropropane	ND	0.400	0.120	ug/L	1	VMS10242	VXX18954
1,2-Dibromoethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Carbon tetrachloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,1,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
Chloroform	ND	1.00	0.300	ug/L	1	VMS10242	VXX18954
Bromobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,3-Trichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromomethane	ND	3.00	0.940	ug/L	1	VMS10242	VXX18954
Bromochloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Vinyl chloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dichlorodifluoromethane	ND	1.00	0.310	ug/L ug/L	1	VMS10242	VXX18954
Distribution of the that is	110	1.00	0.010	ug/L	1	VIVIO IUZ4Z	7,001000-

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Client Sample ID: RHMWA01-WG13

SGS Ref. #: 1085813006

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:05 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

0 1 7	,					<u>Analytical</u>	Prep	
<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	<u>Batch</u>	Batch Qualifie	<u>rs</u>
Chloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
sec-Butylbenzene	8.57	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Bromodichloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954	
1,1-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
2-Butanone (MEK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954	
Methylene chloride	ND	5.00	1.00	ug/L	1	VMS10242	VXX18954	
Trichlorofluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
P & M -Xylene	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954	
Naphthalene	245	40.0	12.4	ug/L	20	VMS10254	VXX18976	
o-Xylene	0.490 J	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Bromoform	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1-Chlorohexane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
tert-Butylbenzene	1.31	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1-Dichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
2-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Trichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
2,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Hexachlorobutadiene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Isopropylbenzene (Cumene)	6.48	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2-Dichloroethane-D4 <surr></surr>	103	73-120		%	1	VMS10242	VXX18954	
Toluene-d8 <surr></surr>	103	80-120		%	1	VMS10242	VXX18954	
4-Bromofluorobenzene <surr></surr>	105	76-120		%	1	VMS10242	VXX18954	



Client Sample ID: RHMWA01-WG13

SGS Ref. #: 1085813006

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:05 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

						<u>Analytical</u>	Prep	
<u>Parameter</u>	<u>Result</u>	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	<u>Batch</u>	<u>Batch</u>	<b>Qualifiers</b>
Batch Information								
Analytical Batch: VMS10242		Prep Batch	ı: VXX18954			Initial Prep	Wt./Vol.: 5	mL
Analytical Method: SW8260B		Prep Metho	od: SW5030B			Prep Extrac	t Vol.: 5 m	L
Analysis Date/Time: 10/27/08 18:58		Prep Date/	Time: 10/27/08	11:41		Container II	D:1085813	006-D
Dilution Factor: 1						Analyst: JD	В	
Analytical Batch: VMS10254		Prep Batch	ı: VXX18976			Initial Prep	Wt./Vol.: 5	mL
Analytical Method: SW8260B		Prep Metho	od: SW5030B			Prep Extrac	t Vol.: 5 m	L
Analysis Date/Time: 10/31/08 09:09		Prep Date/	Time: 10/31/08	01:28		Container II	D:1085813	006-C
Dilution Factor: 20						Analyst: DS	SH	



Client Sample ID: RHMWA01-WG13

SGS Ref. #: 1085813006

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 12:05 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### Polynuclear Aromatics GC/MS

<u>Parameter</u>	<u>Result</u>	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Analytical Batch	Prep Batch Qualifiers
Acenaphthylene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Acenaphthene	0.208	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Fluorene	0.122	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Phenanthrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Anthracene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Fluoranthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Pyrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo(a)Anthracene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Chrysene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[b]Fluoranthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[k]fluoranthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[a]pyrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Indeno[1,2,3-c,d] pyrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Dibenzo[a,h]anthracene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[g,h,i]perylene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Naphthalene	82.3	5.00	1.55	ug/L	50	XMS4763	XXX20276
1-Methylnaphthalene	62.4	2.50	0.750	ug/L	50	XMS4763	XXX20276
2-Methylnaphthalene	12.7	1.00	0.300	ug/L	20	XMS4763	XXX20276
Terphenyl-d14 <surr></surr>	87.7	50-135		%	1	XMS4761	XXX20276
Batch Information							
Analytical Batch: XMS4761		•	: XXX20276				Wt./Vol.: 1000 mL
Analytical Method: 8270D SIMS		•	od: SW3520C	0.40		•	et Vol.: 1 mL
Analysis Date/Time: 10/30/08 17:36 Dilution Factor: 1		Prep Date/	Time: 10/28/08 1	0:10		Analyst: JD	D:1085813006-J iH
Analytical Batch: XMS4763		Prep Batch	: XXX20276			Initial Prep	Wt./Vol.: 1000 mL
Analytical Method: 8270D SIMS		Prep Metho	od: SW3520C			Prep Extrac	t Vol.: 1 mL
Analysis Date/Time: 10/31/08 12:22		Prep Date/	Time: 10/28/08 1	0:10		Container II	D:1085813006-J
Dilution Factor: 20						Analyst: JD	H
Analytical Batch: XMS4763		Prep Batch	: XXX20276			Initial Prep	Wt./Vol.: 1000 mL
Analytical Method: 8270D SIMS		Prep Metho	od: SW3520C			Prep Extrac	ct Vol.: 1 mL
Analysis Date/Time: 10/31/08 16:31		Prep Date/	Time: 10/28/08 1	0:10		Container II	D:1085813006-J
Dilution Factor: 50						Analyst: JD	H



Client Sample ID: RHMW01-WG13

SGS Ref. #: 1085813007

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 15:00 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Fuels Department**

<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> Batch	Qualifiers
Gasoline Range Organics	ND	100	10.0	ug/L	1	VFC9248	VXX1896	7
4-Bromofluorobenzene <surr></surr>	92.5	50-150		%	1	VFC9248	VXX1896	7
Batch Information								
Analytical Batch: VFC9248		Prep Batch	: VXX18967			Initial Prep	Wt./Vol.: 5 n	nL
Analytical Method: SW8015C		Prep Metho	od: SW5030B			Prep Extra	ct Vol.: 5 mL	
Analysis Date/Time: 10/30/08 12:12		Prep Date/	Time: 10/30/08	08:00		Container I	D:10858130	007-A
Dilution Factor: 1						Analyst: HI	M	



Client Sample ID: RHMW01-WG13

SGS Ref. #: 1085813007

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 15:00 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

Analytical

Prep

### **Semivolatile Organic Fuels Department**

<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Batch	Batch	<u>Qualifiers</u>
Diesel Range Organics	0.459	0.412	0.0825	mg/L	1	XFC8323	XXX2029	0
5a Androstane <surr></surr>	103	50-150		%	1	XFC8323	XXX2029	0
Batch Information								
Analytical Batch: XFC8323		Prep Batch	: XXX20290			Initial Prep	Wt./Vol.: 970	0 mL
Analytical Method: SW8015C		Prep Metho	od: SW3520C			Prep Extra	ct Vol.: 1 mL	
Analysis Date/Time: 10/31/08 13:01		Prep Date/	Time: 10/29/08 1	8:00		Container I	D:10858130	07-H
Dilution Factor: 1						Analyst: G	L	



Client Sample ID: RHMW01-WG13

SGS Ref. #: 1085813007

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 15:00 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

						Analytical	Prep
<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Benzene	ND	0.400	0.120	ug/L	1	VMS10242	VXX18954
Toluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Ethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
n-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dichloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Isopropyltoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
n-Propylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Styrene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromomethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,4-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Acetone	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dibromo-3-chloropropane	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
Methyl-t-butyl ether	ND	5.00	1.50	ug/L	1	VMS10242	VXX18954
Tetrachloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromochloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3-Dichloropropane	ND	0.400	0.120	ug/L	1	VMS10242	VXX18954
1,2-Dibromoethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Carbon tetrachloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,1,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
Chloroform	ND	1.00	0.300	ug/L	1	VMS10242	VXX18954
Bromobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,3-Trichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromomethane	ND	3.00	0.940	ug/L	1	VMS10242	VXX18954
Bromochloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Vinyl chloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dichlorodifluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954

SGS Environmental Services Inc. Alaska Division 200 West Potter Drive Anchorage Alaska 99518

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Client Sample ID: RHMW01-WG13

SGS Ref. #: 1085813007

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 15:00 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

Parameter	Result	PQL/CL	MDL	Units	<u>DF</u>	Analytical Batch	Prep Batch Qualifiers
Chloroethane	ND	1.00	0.310		<u> </u>	VMS10242	VXX18954
sec-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242 VMS10242	VXX18954 VXX18954
Bromodichloromethane	ND	0.500	0.150	ug/L ug/L	1	VMS10242 VMS10242	VXX18954 VXX18954
1,1-Dichloroethene	ND	1.00	0.130	_		VMS10242 VMS10242	VXX18954 VXX18954
•	ND	10.0	3.10	ug/L	1		VXX18954 VXX18954
2-Butanone (MEK)	ND ND	5.00	1.00	ug/L	1	VMS10242	VXX18954 VXX18954
Methylene chloride				ug/L	1	VMS10242	
Trichlorofluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
P & M -Xylene	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
Naphthalene	ND	2.00	0.620	ug/L	1	VMS10254	VXX18976
o-Xylene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromoform	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1-Chlorohexane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
tert-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1-Dichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
2-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Trichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
2,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Hexachlorobutadiene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Isopropylbenzene (Cumene)	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2-Dichloroethane-D4 <surr></surr>	103	73-120	0.0.0	%	1	VMS10242	VXX18954
Toluene-d8 <surr></surr>	104	80-120		%	1	VMS10242 VMS10242	VXX18954
4-Bromofluorobenzene <surr></surr>	109	76-120		%	1	VMS10242 VMS10242	VXX18954



Client Sample ID: RHMW01-WG13

SGS Ref. #: 1085813007

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 15:00 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

Parameter	Result	PQL/CL	MDL_	Units	DF	Analytical Batch	Prep Batch	Qualifiers		
Batch Information					_					
Analytical Batch: VMS10242		Prep Batch	ı: VXX18954			Initial Prep	Wt./Vol.: 5	mL		
Analytical Method: SW8260B	Prep Method: SW5030B					Prep Extract Vol.: 5 mL				
Analysis Date/Time: 10/27/08 19:31		Prep Date/Time: 10/27/08 11:41				Container ID:1085813007-D				
Dilution Factor: 1						Analyst: JD	В			
Analytical Batch: VMS10254		Prep Batch	ı: VXX18976			Initial Prep	Wt./Vol.: 5	mL		
Analytical Method: SW8260B		Prep Method: SW5030B				Prep Extract Vol.: 5 mL				
Analysis Date/Time: 10/31/08 07:36		Prep Date/	Time: 10/31/08	01:28		Container II	D:1085813	007-F		
Dilution Factor: 1						Analyst: DS	SH			



Client Sample ID: RHMW01-WG13

SGS Ref. #: 1085813007

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 15:00 Receipt Date/Time: 10/24/08 10:50

### Polynuclear Aromatics GC/MS

<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Acenaphthylene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Acenaphthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Fluorene	0.0207 J	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Phenanthrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Anthracene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Fluoranthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Pyrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo(a)Anthracene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Chrysene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[b]Fluoranthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[k]fluoranthene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[a]pyrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Indeno[1,2,3-c,d] pyrene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Dibenzo[a,h]anthracene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Benzo[g,h,i]perylene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Naphthalene	0.103	0.100	0.0310	ug/L	1	XMS4761	XXX20276
1-Methylnaphthalene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
2-Methylnaphthalene	ND	0.0500	0.0150	ug/L	1	XMS4761	XXX20276
Terphenyl-d14 <surr></surr>	91.1	50-135		%	1	XMS4761	XXX20276

#### **Batch Information**

Prep Batch: XXX20276 Initial Prep Wt./Vol.: 1000 mL Analytical Batch: XMS4761 Analytical Method: 8270D SIMS Prep Method: SW3520C Prep Extract Vol.: 1 mL Analysis Date/Time: 10/30/08 18:09 Prep Date/Time: 10/28/08 10:10 Container ID:1085813007-J Dilution Factor: 1

Analyst: JDH

Analytical

Prep

Print Date: 11/20/2008



Client Sample ID: TB01-WG13

SGS Ref. #: 1085813008

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 10/22/08 08:05 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

### **Volatile Gas Chromatography/Mass Spectroscopy**

, , , , , , , , , , , , , , , , , , ,						Analytical	Prep
<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Benzene	ND	0.400	0.120	ug/L	1	VMS10242	VXX18954
Toluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Ethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
n-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dichloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chlorobenzene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
4-Isopropyltoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
n-Propylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Styrene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromomethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,4-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Acetone	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954
1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,2-Dibromo-3-chloropropane	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954
Methyl-t-butyl ether	ND	5.00	1.50	ug/L	1	VMS10242	VXX18954
Tetrachloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dibromochloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
1,3-Dichloropropane	ND	0.400	0.120	ug/L	1	VMS10242	VXX18954
1,2-Dibromoethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Carbon tetrachloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,1,1,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954
Chloroform	ND	1.00	0.300	ug/L	1	VMS10242	VXX18954
Bromobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Chloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
1,2,3-Trichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Bromomethane	ND	3.00	0.940	ug/L	1	VMS10242	VXX18954
Bromochloromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Vinyl chloride	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954
Dichlorodifluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954

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Client Sample ID: TB01-WG13 SGS Ref. #: 1085813008

Project ID: 9121-003 Red Hill BFSF

Matrix: Water (Surface, Eff., Ground)

Collection Date/Time: 10/22/08 08:05 Receipt Date/Time: 10/24/08 10:50

Print Date: 11/20/2008

Analytical

Prep

### Volatile Gas Chromatography/Mass Spectroscopy

						Analytical	<u>Prep</u>	
<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	DF	<u>Batch</u>		<u>Qualifiers</u>
Chloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
sec-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Bromodichloromethane	ND	0.500	0.150	ug/L	1	VMS10242	VXX18954	
1,1-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
2-Butanone (MEK)	ND	10.0	3.10	ug/L	1	VMS10242	VXX18954	
Methylene chloride	ND	5.00	1.00	ug/L	1	VMS10242	VXX18954	
Trichlorofluoromethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
P & M -Xylene	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954	
Naphthalene	ND	2.00	0.620	ug/L	1	VMS10242	VXX18954	
o-Xylene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Bromoform	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1-Chlorohexane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
tert-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1-Dichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
2-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Trichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
2,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Hexachlorobutadiene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
Isopropylbenzene (Cumene)	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10242	VXX18954	
1,2-Dichloroethane-D4 <surr></surr>	105	73-120		%	1	VMS10242	VXX18954	
Toluene-d8 <surr></surr>	106	80-120		%	1	VMS10242	VXX18954	
4-Bromofluorobenzene <surr></surr>	112	76-120		%	1	VMS10242	VXX18954	

#### **Batch Information**

Prep Batch: VXX18954 Initial Prep Wt./Vol.: 5 mL Analytical Batch: VMS10242 Analytical Method: SW8260B Prep Method: SW5030B Prep Extract Vol.: 5 mL Analysis Date/Time: 10/27/08 16:09 Prep Date/Time: 10/27/08 11:41 Container ID:1085813008-A

Dilution Factor: 1 Analyst: JDB



SGS Ref.#

867629

Method Blank

Client Name

The Environmental Company, Inc. (TEC)

Project Name/# Matrix 9121-003 Red Hill BFSF

Water (Surface, Eff., Ground)

Printed Date/Time
Prep Batch

11/20/2008 8:09

Method

VXX18954 SW5030B

Date 10/27/2008

QC results affect the following production samples:

1085813001, 1085813004, 1085813005, 1085813006, 1085813007, 1085813008

Politic Gas Chromatography/Mass   Spectroscopy	Parameter	Results	Reporting/Control Limit	MDL	Units	Analysis Date
Toluene ND 1.00 0.310 ug/L 102708 Ethylbenzene ND 1.00 0.310 ug/L 102708 I-Butylbenzene ND 1.00 0.310 ug/L 102708 I-Butylbenzene ND 0.500 0.150 ug/L 102708 I-J-Dichloroetnane ND 0.500 0.150 ug/L 102708 I-J-Dichloroetnane ND 1.00 0.310 ug/L 102708 I-J-S-Trimethylbenzene ND 1.00 0.310 ug/L 102708 I-J-Dichloroethane ND 1.00 0.310 ug/L 102708 I-J-Dichloroptopene ND 1.00 0.310 ug/L 102708 I-J-Z-Tritachloroethane ND 0.500 0.150 ug/L 102708 I-J-Dichloroptopane ND 0.500 0.150 ug/L 102708 I-J-Dichl	Volatile Gas Chromatography/	Mass Spectr	oscopy			
Hithylbenzene   ND   1.00   0.310   ug/L   1027/08	Benzene	ND	0.400	0.120	ug/L	10/27/08
n-Butylbenzene         ND         1.00         0.310         ug/L         10.27/08           1,4-Dichlorobenzene         ND         0.500         0.150         ug/L         10.27/08           1,2-Dichloroethane         ND         0.500         0.150         ug/L         10.27/08           1,3-5-Trimethylbenzene         ND         1.00         0.310         ug/L         10.27/08           4-Chlorotoluene         ND         1.00         0.310         ug/L         10.27/08           4-Methyl-2-pentanone (MIBK)         ND         1.00         3.10         ug/L         10.27/08           4-Methyl-2-pentanone (MIBK)         ND         1.00         0.310         ug/L         10.27/08           4-Hothyl-2-pentanone (MIBK)         ND         1.00         0.310         ug/L         10.27/08           4-Isopropylloluene         ND         1.00         0.310         ug/L         10.27/08           4-Isopropylloluene         ND         1.00         0.310         ug/L         10.27/08           6is-1,3-Dichloropropene         ND         1.00         0.310         ug/L         10.27/08           8tyrene         ND         1.00         0.310         ug/L         10.27/08	Toluene	ND	1.00	0.310	ug/L	10/27/08
1,4-Dichlorobenzene	Ethylbenzene	ND	1.00	0.310	ug/L	10/27/08
1,2-Dichloroethane	n-Butylbenzene	ND	1.00	0.310	ug/L	10/27/08
1,3,5-Tirmethylbenzene         ND         1.00         0.310         ug/L         1027/08           4-Chlorotoluene         ND         1.00         0.310         ug/L         1027/08           Chlorobenzene         ND         0.500         0.150         ug/L         1027/08           4-Methyl-2-pentanon (MIBK)         ND         1.00         0.310         ug/L         1027/08           6is-1,2-Dichlorochene         ND         1.00         0.310         ug/L         1027/08           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1027/08           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1027/08           8-Propylbenzene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Lay-Tirchloropropene         ND         1.00         0.310         ug/L         1027/08           Acetone         ND	1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	10/27/08
4-Chlorotoluene         ND         1.00         0.310         ug/L         1027/08           Chlorobenzene         ND         0.500         0.150         ug/L         1027/08           4-Methyl-2-pentanone (MIBK)         ND         1.00         3.10         ug/L         1027/08           4-Methyl-2-pentanone (MIBK)         ND         1.00         3.10         ug/L         1027/08           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1027/08           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1027/08           cis-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           strans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1027/08           1,2,2-Tircheloropropene         ND         1.00         0.310         ug/L         1027/08           Acetone	1,2-Dichloroethane	ND	0.500	0.150	ug/L	10/27/08
Chlorobenzene         ND         0.500         0.150         ug/L         1027/08           4-Methyl-2-pentanone (MIBK)         ND         10.0         3.10         ug/L         1027/08           4-Methyl-2-pentanone (MIBK)         ND         1.00         0.310         ug/L         1027/08           4-Isopropytholuene         ND         1.00         0.310         ug/L         1027/08           6-Isopropytholuene         ND         1.00         0.310         ug/L         1027/08           6-Isopropytholuene         ND         1.00         0.310         ug/L         1027/08           6-Propythenzene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1027/08           1,2,2-Tetrachloroethane         ND         1.00         0.310         ug/L         1027/08           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1027/08           Methyl-t-buty	1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	10/27/08
4-Methyl-2-pentanone (MIBK)         ND         10.0         3.10         ug/L         1027/08           cis-1,2-Dichloroethene         ND         1.00         0.310         ug/L         1027/08           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1027/08           cis-1,3-Dichloropropene         ND         0.500         0.150         ug/L         1027/08           n-Propylbenzene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1027/08           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1027/08           4-Ectone         ND         1.00         3.10         ug/L         1027/08           1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1027/08           1,2-Dibromo-3-chloropropane         ND         5.00         1.50         ug/L         1027/08           Methyl-1-but	4-Chlorotoluene	ND	1.00	0.310	ug/L	10/27/08
cis-1,2-Dichloroethene         ND         1.00         0.310         ug/L         1027/08           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1027/08           cis-1,3-Dichloropropene         ND         0.500         0.150         ug/L         1027/08           N-Propylbenzene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           bibromonethane         ND         1.00         0.310         ug/L         1027/08           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1027/08           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1027/08           1,2,2-Tetrachloroethane         ND         1.00         3.10         ug/L         1027/08           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1027/08           1,2-Dibromo-3-chloropropane         ND         5.00         1.50         ug/L         1027/08           1,2-Dibromo-3-chloropropane         ND         0.500         0.150         ug/L         1027/08 <t< td=""><td>Chlorobenzene</td><td>ND</td><td>0.500</td><td>0.150</td><td>ug/L</td><td>10/27/08</td></t<>	Chlorobenzene	ND	0.500	0.150	ug/L	10/27/08
cis-1,2-Dichloroethene         ND         1.00         0.310         ug/L         1027/08           4-Isopropyltoluene         ND         1.00         0.310         ug/L         1027/08           cis-1,3-Dichloropropene         ND         0.500         0.150         ug/L         1027/08           Propylbenzene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Dibromomethane         ND         1.00         0.310         ug/L         1027/08           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1027/08           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1027/08           1,2,2-Tetrachloroethane         ND         1.00         0.310         ug/L         1027/08           1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1027/08           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1027/08           Tetrachloroethane         ND         0.500         0.150         ug/L         1027/08           1,3-Di	4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	10/27/08
cis-1,3-Dichloropropene         ND         0.500         0.150         ug/L         1027/08           n-Propylbenzene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Dibromomethane         ND         1.00         0.310         ug/L         1027/08           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1027/08           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1027/08           Acetone         ND         1.00         3.10         ug/L         1027/08           1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1027/08           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1027/08           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1027/08           Methyl-t-butyl ether         ND         0.500         0.150         ug/L         1027/08           Tetrachloroethane         ND         0.500         0.150         ug/L         1027/08           1,2-Dibromoch		ND	1.00	0.310	ug/L	10/27/08
cis-1,3-Dichloropropene         ND         0.500         0.150         ug/L         1027/08           n-Propylbenzene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Dibromomethane         ND         1.00         0.310         ug/L         1027/08           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1027/08           L2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1027/08           Acetone         ND         10.0         3.10         ug/L         1027/08           1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1027/08           1,2,2-Dibromo-3-chloropropane         ND         0.500         0.150         ug/L         1027/08           Methyl-t-butyl ether         ND         0.500         0.150         ug/L         1027/08           Tetrachloroethane         ND         0.500         0.150         ug/L         1027/08           1,3-Dichloropropane         ND         0.400         0.310         ug/L         1027/08           1,2-Dibrom	4-Isopropyltoluene	ND	1.00	0.310	ug/L	10/27/08
n-Propylbenzene         ND         1.00         0.310         ug/L         1027/08           Styrene         ND         1.00         0.310         ug/L         1027/08           Dibromomethane         ND         1.00         0.310         ug/L         1027/08           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         1027/08           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         1027/08           Acetone         ND         1.00         3.10         ug/L         1027/08           1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         1027/08           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         1027/08           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         1027/08           Methyl-t-butyl ether         ND         1.00         0.310         ug/L         1027/08           Tetrachloroethane         ND         1.00         0.310         ug/L         1027/08           1,3-Dichloropropane         ND         0.400         0.120         ug/L         1027/08           1,1,1,2-Tetrachloro		ND	0.500	0.150	ug/L	10/27/08
Styrene         ND         1.00         0.310         ug/L         1027/08           Dibromomethane         ND         1.00         0.310         ug/L         1027/08           trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         10/27/08           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         10/27/08           Acetone         ND         10.0         3.10         ug/L         10/27/08           1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         10/27/08           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         10/27/08           Tetrachloroethene         ND         1.00         0.310         ug/L         10/27/08           Dibromochloromethane         ND         0.500         0.150         ug/L         10/27/08           1,2-Dibromochane         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         1.00         0.310         ug/L         10/27/08		ND	1.00	0.310	ug/L	10/27/08
trans-1,3-Dichloropropene         ND         1.00         0.310         ug/L         10/27/08           1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         10/27/08           Acetone         ND         10.0         3.10         ug/L         10/27/08           1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         10/27/08           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         10/27/08           Tetrachloroethene         ND         1.00         0.310         ug/L         10/27/08           Tetrachloroethene         ND         0.500         0.150         ug/L         10/27/08           1,3-Dichloropropane         ND         0.400         0.120         ug/L         10/27/08           1,2-Dibromoethane         ND         1.00         0.310         ug/L         10/27/08           Carbon tetrachloride         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         1.00         0.310         ug/L         10/27/08		ND	1.00	0.310	ug/L	10/27/08
1,2,4-Trichlorobenzene	Dibromomethane	ND	1.00	0.310	ug/L	10/27/08
1,2,4-Trichlorobenzene         ND         1.00         0.310         ug/L         10/27/08           Acetone         ND         10.0         3.10         ug/L         10/27/08           1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         10/27/08           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         10/27/08           Tetrachloroethene         ND         1.00         0.310         ug/L         10/27/08           Dibromochloromethane         ND         0.500         0.150         ug/L         10/27/08           1,3-Dichloropropane         ND         0.400         0.120         ug/L         10/27/08           1,2-Dibromoethane         ND         1.00         0.310         ug/L         10/27/08           Carbon tetrachloride         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         1.00         0.300         ug/L         10/27/08           Bromobenzene         ND         1.00         0.310         ug/L         10/27/08	trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	10/27/08
Acetone         ND         10.0         3.10         ug/L         10/27/08           1,1,2,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         10/27/08           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         10/27/08           Tetrachloroethene         ND         1.00         0.310         ug/L         10/27/08           Dibromochloromethane         ND         0.500         0.150         ug/L         10/27/08           1,3-Dichloropropane         ND         0.400         0.120         ug/L         10/27/08           1,2-Dibromoethane         ND         1.00         0.310         ug/L         10/27/08           1,2-Dibromoethane         ND         1.00         0.310         ug/L         10/27/08           Carbon tetrachloride         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         1.00         0.310         ug/L         10/27/08           Chloroform         ND         1.00         0.310         ug/L         10/27/08 <t< td=""><td></td><td>ND</td><td>1.00</td><td>0.310</td><td>ug/L</td><td>10/27/08</td></t<>		ND	1.00	0.310	ug/L	10/27/08
1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         10/27/08           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         10/27/08           Tetrachloroethene         ND         1.00         0.310         ug/L         10/27/08           Dibromochloromethane         ND         0.500         0.150         ug/L         10/27/08           1,3-Dichloropropane         ND         0.400         0.120         ug/L         10/27/08           1,2-Dibromoethane         ND         1.00         0.310         ug/L         10/27/08           Carbon tetrachloride         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           Chloroform         ND         1.00         0.300         ug/L         10/27/08           Bromobenzene         ND         1.00         0.310         ug/L         10/27/08           Chloromethane         ND         1.00         0.310         ug/L         10/27/08           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         10/27/08 <td< td=""><td></td><td>ND</td><td>10.0</td><td>3.10</td><td>ug/L</td><td>10/27/08</td></td<>		ND	10.0	3.10	ug/L	10/27/08
1,2-Dibromo-3-chloropropane         ND         2.00         0.620         ug/L         10/27/08           Methyl-t-butyl ether         ND         5.00         1.50         ug/L         10/27/08           Tetrachloroethene         ND         1.00         0.310         ug/L         10/27/08           Dibromochloromethane         ND         0.500         0.150         ug/L         10/27/08           1,3-Dichloropropane         ND         0.400         0.120         ug/L         10/27/08           1,2-Dibromoethane         ND         1.00         0.310         ug/L         10/27/08           Carbon tetrachloride         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           Chloroform         ND         1.00         0.300         ug/L         10/27/08           Bromobenzene         ND         1.00         0.310         ug/L         10/27/08           Chloromethane         ND         1.00         0.310         ug/L         10/27/08           1,2,3-Trichloropropane         ND         3.00         0.940         ug/L         10/27/08 <td< td=""><td>1,1,2,2-Tetrachloroethane</td><td>ND</td><td>0.500</td><td>0.150</td><td>ug/L</td><td>10/27/08</td></td<>	1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	10/27/08
Methyl-t-butyl ether         ND         5.00         1.50         ug/L         10/27/08           Tetrachloroethene         ND         1.00         0.310         ug/L         10/27/08           Dibromochloromethane         ND         0.500         0.150         ug/L         10/27/08           1,3-Dichloropropane         ND         0.400         0.120         ug/L         10/27/08           1,2-Dibromoethane         ND         1.00         0.310         ug/L         10/27/08           Carbon tetrachloride         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           Chloroform         ND         1.00         0.300         ug/L         10/27/08           Bromobenzene         ND         1.00         0.310         ug/L         10/27/08           Chloromethane         ND         1.00         0.310         ug/L         10/27/08           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         10/27/08           Bromochloromethane         ND         1.00         0.310         ug/L         10/27/08           Vinyl ch		ND	2.00	0.620	ug/L	10/27/08
Tetrachloroethene         ND         1.00         0.310         ug/L         10/27/08           Dibromochloromethane         ND         0.500         0.150         ug/L         10/27/08           1,3-Dichloropropane         ND         0.400         0.120         ug/L         10/27/08           1,2-Dibromoethane         ND         1.00         0.310         ug/L         10/27/08           Carbon tetrachloride         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           Chloroform         ND         1.00         0.300         ug/L         10/27/08           Bromobenzene         ND         1.00         0.310         ug/L         10/27/08           Chloromethane         ND         1.00         0.310         ug/L         10/27/08           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         10/27/08           Bromochloromethane         ND         3.00         0.940         ug/L         10/27/08           Vinyl chloride         ND         1.00         0.310         ug/L         10/27/08	• •	ND	5.00	1.50	ug/L	10/27/08
Dibromochloromethane         ND         0.500         0.150         ug/L         10/27/08           1,3-Dichloropropane         ND         0.400         0.120         ug/L         10/27/08           1,2-Dibromoethane         ND         1.00         0.310         ug/L         10/27/08           Carbon tetrachloride         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           Chloroform         ND         1.00         0.300         ug/L         10/27/08           Bromobenzene         ND         1.00         0.310         ug/L         10/27/08           Chloromethane         ND         1.00         0.310         ug/L         10/27/08           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         10/27/08           Bromochloromethane         ND         3.00         0.940         ug/L         10/27/08           Vinyl chloride         ND         1.00         0.310         ug/L         10/27/08	-	ND	1.00	0.310		10/27/08
1,3-Dichloropropane       ND       0.400       0.120       ug/L       10/27/08         1,2-Dibromoethane       ND       1.00       0.310       ug/L       10/27/08         Carbon tetrachloride       ND       1.00       0.310       ug/L       10/27/08         1,1,1,2-Tetrachloroethane       ND       0.500       0.150       ug/L       10/27/08         Chloroform       ND       1.00       0.300       ug/L       10/27/08         Bromobenzene       ND       1.00       0.310       ug/L       10/27/08         Chloromethane       ND       1.00       0.310       ug/L       10/27/08         1,2,3-Trichloropropane       ND       1.00       0.310       ug/L       10/27/08         Bromoethane       ND       3.00       0.940       ug/L       10/27/08         Bromochloromethane       ND       1.00       0.310       ug/L       10/27/08         Vinyl chloride       ND       1.00       0.310       ug/L       10/27/08		ND	0.500	0.150	ug/L	10/27/08
1,2-Dibromoethane       ND       1.00       0.310       ug/L       10/27/08         Carbon tetrachloride       ND       1.00       0.310       ug/L       10/27/08         1,1,1,2-Tetrachloroethane       ND       0.500       0.150       ug/L       10/27/08         Chloroform       ND       1.00       0.300       ug/L       10/27/08         Bromobenzene       ND       1.00       0.310       ug/L       10/27/08         Chloromethane       ND       1.00       0.310       ug/L       10/27/08         1,2,3-Trichloropropane       ND       1.00       0.310       ug/L       10/27/08         Bromoethane       ND       3.00       0.940       ug/L       10/27/08         Bromochloromethane       ND       1.00       0.310       ug/L       10/27/08         Vinyl chloride       ND       1.00       0.310       ug/L       10/27/08	1,3-Dichloropropane	ND	0.400	0.120		10/27/08
Carbon tetrachloride         ND         1.00         0.310         ug/L         10/27/08           1,1,1,2-Tetrachloroethane         ND         0.500         0.150         ug/L         10/27/08           Chloroform         ND         1.00         0.300         ug/L         10/27/08           Bromobenzene         ND         1.00         0.310         ug/L         10/27/08           Chloromethane         ND         1.00         0.310         ug/L         10/27/08           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         10/27/08           Bromoethane         ND         3.00         0.940         ug/L         10/27/08           Bromochloromethane         ND         1.00         0.310         ug/L         10/27/08           Vinyl chloride         ND         1.00         0.310         ug/L         10/27/08		ND	1.00	0.310		10/27/08
1,1,1,2-Tetrachloroethane       ND       0.500       0.150       ug/L       10/27/08         Chloroform       ND       1.00       0.300       ug/L       10/27/08         Bromobenzene       ND       1.00       0.310       ug/L       10/27/08         Chloromethane       ND       1.00       0.310       ug/L       10/27/08         1,2,3-Trichloropropane       ND       1.00       0.310       ug/L       10/27/08         Bromomethane       ND       3.00       0.940       ug/L       10/27/08         Bromochloromethane       ND       1.00       0.310       ug/L       10/27/08         Vinyl chloride       ND       1.00       0.310       ug/L       10/27/08	*	ND	1.00			10/27/08
Chloroform         ND         1.00         0.300         ug/L         10/27/08           Bromobenzene         ND         1.00         0.310         ug/L         10/27/08           Chloromethane         ND         1.00         0.310         ug/L         10/27/08           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         10/27/08           Bromomethane         ND         3.00         0.940         ug/L         10/27/08           Bromochloromethane         ND         1.00         0.310         ug/L         10/27/08           Vinyl chloride         ND         1.00         0.310         ug/L         10/27/08		ND	0.500	0.150		10/27/08
Bromobenzene         ND         1.00         0.310         ug/L         10/27/08           Chloromethane         ND         1.00         0.310         ug/L         10/27/08           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         10/27/08           Bromomethane         ND         3.00         0.940         ug/L         10/27/08           Bromochloromethane         ND         1.00         0.310         ug/L         10/27/08           Vinyl chloride         ND         1.00         0.310         ug/L         10/27/08		ND	1.00			10/27/08
Chloromethane         ND         1.00         0.310         ug/L         10/27/08           1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         10/27/08           Bromomethane         ND         3.00         0.940         ug/L         10/27/08           Bromochloromethane         ND         1.00         0.310         ug/L         10/27/08           Vinyl chloride         ND         1.00         0.310         ug/L         10/27/08			1.00			10/27/08
1,2,3-Trichloropropane         ND         1.00         0.310         ug/L         10/27/08           Bromomethane         ND         3.00         0.940         ug/L         10/27/08           Bromochloromethane         ND         1.00         0.310         ug/L         10/27/08           Vinyl chloride         ND         1.00         0.310         ug/L         10/27/08						10/27/08
Bromomethane         ND         3.00         0.940         ug/L         10/27/08           Bromochloromethane         ND         1.00         0.310         ug/L         10/27/08           Vinyl chloride         ND         1.00         0.310         ug/L         10/27/08			1.00			10/27/08
Bromochloromethane         ND         1.00         0.310         ug/L         10/27/08           Vinyl chloride         ND         1.00         0.310         ug/L         10/27/08			3.00			10/27/08
Vinyl chloride ND 1.00 0.310 ug/L 10/27/08			1.00		=	10/27/08
			1.00			10/27/08
			1.00			10/27/08



SGS Ref.# Client Name 867629 Method Blank

The Environmental Company, Inc. (TEC)

Project Name/# Matrix 9121-003 Red Hill BFSF Water (Surface, Eff., Ground)

VMS10242

SW8260B

HP 5890 Series II MS3 VNA

Batch Method

Instrument

Printed Date/Time
Prep Batch

11/20/2008 8:09 VXX18954

Method Date

SW5030B 10/27/2008

Parameter	Results	Reporting/Control	MDL	Units	Analysis Date
Volatile Gas Chromatography/M	lass Spectro	saony			
VOIACITE GAS CHIOMACOGIAPHY/F	ass spectio	scopy			
Chloroethane	ND	1.00	0.310	ug/L	10/27/08
sec-Butylbenzene	ND	1.00	0.310	ug/L	10/27/08
Bromodichloromethane	ND	0.500	0.150	ug/L	10/27/08
1,1-Dichloroethene	ND	1.00	0.310	ug/L	10/27/08
2-Butanone (MEK)	ND	10.0	3.10	ug/L	10/27/08
Methylene chloride	ND	5.00	1.00	ug/L	10/27/08
Trichlorofluoromethane	ND	1.00	0.310	ug/L	10/27/08
P & M -Xylene	ND	2.00	0.620	ug/L	10/27/08
Naphthalene	ND	2.00	0.620	ug/L	10/27/08
o-Xylene	ND	1.00	0.310	ug/L	10/27/08
Bromoform	ND	1.00	0.310	ug/L	10/27/08
1-Chlorohexane	ND	1.00	0.310	ug/L	10/27/08
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	10/27/08
tert-Butylbenzene	ND	1.00	0.310	ug/L	10/27/08
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	10/27/08
1,1-Dichloroethane	ND	1.00	0.310	ug/L	10/27/08
2-Chlorotoluene	ND	1.00	0.310	ug/L	10/27/08
Trichloroethene	ND	1.00	0.310	ug/L	10/27/08
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	10/27/08
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	10/27/08
2,2-Dichloropropane	ND	1.00	0.310	ug/L	10/27/08
Hexachlorobutadiene	ND	1.00	0.310	ug/L	10/27/08
Isopropylbenzene (Cumene)	ND	1.00	0.310	ug/L	10/27/08
1,2-Dichloropropane	ND	1.00	0.310	ug/L	10/27/08
1,1-Dichloropropene	ND	1.00	0.310	ug/L	10/27/08
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	10/27/08
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	10/27/08
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	10/27/08
Surrogates					
1,2-Dichloroethane-D4 <surr></surr>	102	73-120		%	10/27/08
Toluene-d8 <surr></surr>	105	80-120		%	10/27/08
4-Bromofluorobenzene <surr></surr>	109	76-120		%	10/27/08



SGS Ref.#

867655

Method Blank

The Environmental Company, Inc. (TEC)

**Printed Date/Time** Batch Prep

11/20/2008 8:09

**Client Name** 

9121-003 Red Hill BFSF

Method

Date

XXX20276 SW3520C 10/28/2008

Project Name/# Matrix

Water (Surface, Eff., Ground)

QC results affect the following production samples:

1085813001, 1085813004, 1085813005, 1085813006, 1085813007

Parameter		Limit	MDL	Units	Analysis Date
Polynuclear Aromatics GC/MS					
Acenaphthylene	ND	0.0500	0.0150	ug/L	10/30/08
Acenaphthene	ND	0.0500	0.0150	ug/L	10/30/08
Fluorene	ND	0.0500	0.0150	ug/L	10/30/08
Phenanthrene	ND	0.0500	0.0150	ug/L	10/30/08
Anthracene	ND	0.0500	0.0150	ug/L	10/30/08
Fluoranthene	ND	0.0500	0.0150	ug/L	10/30/08
Pyrene	ND	0.0500	0.0150	ug/L	10/30/08
Benzo(a)Anthracene	ND	0.0500	0.0150	ug/L	10/30/08
Chrysene	ND	0.0500	0.0150	ug/L	10/30/08
Benzo[b]Fluoranthene	ND	0.0500	0.0150	ug/L	10/30/08
Benzo[k]fluoranthene	ND	0.0500	0.0150	ug/L	10/30/08
Benzo[a]pyrene	ND	0.0500	0.0150	ug/L	10/30/08
Indeno[1,2,3-c,d] pyrene	ND	0.0500	0.0150	ug/L	10/30/08
Dibenzo[a,h]anthracene	ND	0.0500	0.0150	ug/L	10/30/08
Benzo[g,h,i]perylene	ND	0.0500	0.0150	ug/L	10/30/08
Naphthalene	ND	0.100	0.0310	ug/L	10/30/08
1-Methylnaphthalene	ND	0.0500	0.0150	ug/L	10/30/08
2-Methylnaphthalene	ND	0.0500	0.0150	ug/L	10/30/08
Surrogates					
Terphenyl-d14 <surr></surr>	88.5	50-135		%	10/30/08

Batch XMS4761 8270D SIMS Method

Instrument HP 5890 Series II MS2 SVOA



SGS Ref.#

868151

Method Blank

Printed Date/Time
Prep Batch

11/20/2008 8:09

Client Name

The Environmental Company, Inc. (TEC)

Batch Method

XXX20290

Project Name/# Matrix 9121-003 Red Hill BFSF

Water (Surface, Eff., Ground)

Date

SW3520C 10/29/2008

QC results affect the following production samples:

1085813004, 1085813005, 1085813006, 1085813007

Parameter		Results	Reporting/Control Limit	MDL	Units	Analysis Date			
Semivolatile Organic Fuels Department									
Diesel Range Org	ganics	ND	0.400	0.0800	mg/L	10/31/08			
Surrogates									
5a Androstane <s< th=""><th>urr&gt;</th><th>104</th><th>60-120</th><th></th><th>%</th><th>10/31/08</th></s<>	urr>	104	60-120		%	10/31/08			
Batch	XFC8323								
Method	SW8015C								
Instrument	HP 6890 Series II FID SV D I	R							



SGS Ref.# Client Name 868324

Method Blank

The Environmental Company, Inc. (TEC)

Project Name/#

9121-003 Red Hill BFSF

Matrix Water (Surface, Eff., Ground)

**Printed Date/Time** 

Prep

11/20/2008 8:09

Batch Method

VXX18967

Date

SW5030B 10/30/2008

QC results affect the following production samples:

1085813001, 1085813004, 1085813005, 1085813006, 1085813007

Parameter		Results	Reporting/Control Limit	MDL	Units	Analysis Date
Volatile Fue	ls Department					
Gasoline Range (	Organics	ND	100	10.0	ug/L	10/30/08
Surrogates						
4-Bromofluorobe	enzene <surr></surr>	94.9	50-150		%	10/30/08
Batch	VFC9248					
Method	SW8015C					
Instrument	HP 5890 Series II PID	+FID VCA				



SGS Ref.#

868746

Method Blank

Printed Date/Time

Prep

11/20/2008 8:09

Client Name

The Environmental Company, Inc. (TEC)

Batch Method VXX18976

Project Name/# Matrix 9121-003 Red Hill BFSF Water (Surface, Eff., Ground)

Date

SW5030B 10/31/2008

QC results affect the following production samples: 1085813005, 1085813006, 1085813007

Parameter		Results	Reporting/Control Limit	MDL	Units	Analysis Date
Volatile Gas	Chromatography/Mass	Spectro	oscopy			
Naphthalene		ND	2.00	0.620	ug/L	10/31/08
Surrogates						
1,2-Dichloroethar	ne-D4 <surr></surr>	102	73-120		%	10/31/08
Toluene-d8 <surr< td=""><td>&gt;</td><td>100</td><td>80-120</td><td></td><td>%</td><td>10/31/08</td></surr<>	>	100	80-120		%	10/31/08
4-Bromofluorobe	nzene <surr></surr>	101	76-120		%	10/31/08
Batch	VMS10254					
Method	SW8260B					
Instrument	HP 5890 Series II MS3 VNA					



SGS Ref.# Client Name 869197

Method Blank

The Environmental Company, Inc. (TEC)

Project Name/# Matrix

9121-003 Red Hill BFSF

Water (Surface, Eff., Ground)

**Printed Date/Time** 

Prep

11/20/2008 8:09

Batch Method

XXX20313 SW3520C

Date 11/03/2008

QC results affect the following production samples:

1085813001

Parameter		Results	Reporting/Control Limit	MDL	Units	Analysis Date
Semivolatile	Organic Fuels Depart	ment				
Diesel Range Org	ganics	ND	0.400	0.0800	mg/L	11/05/08
Surrogates						
5a Androstane <s< td=""><td>urr&gt;</td><td>97.3</td><td>60-120</td><td></td><td>%</td><td>11/05/08</td></s<>	urr>	97.3	60-120		%	11/05/08
Batch	XFC8338					
Method	SW8015C					
Instrument	HD 6900 Sories II FID SV D I	D				

HP 6890 Series II FID SV D R



SGS Ref.#

869581

Method Blank

Printed Date/Time

Prep

11/20/2008 8:09

Client Name

The Environmental Company, Inc. (TEC)

Batch Method

MXX21031 SW3010A

Project Name/# Matrix 9121-003 Red Hill BFSF

Date 11/04/2008

Matrix Water (Surface, Eff., Ground)
QC results affect the following production samples:

1085813001, 1085813004, 1085813005, 1085813006

Reporting/Control Analysis
Parameter Results Limit MDL Units Date

Metals by ICP/MS

Lead ND 1.00 0.310 ug/L 11/14/08

BatchMMS5719MethodSW6020

**Instrument** Perkin Elmer Sciex ICP-MS P4



867630

Lab Control Sample

Printed Date/Time

Prep

11/20/2008

8:09

Client Name

The Environmental Company, Inc. (TEC)

Batch Method VXX18954 SW5030B

Project Name/#
Matrix

9121-003 Red Hill BFSF Water (Surface, Eff., Ground) Date

10/27/2008

QC results affect the following production samples:

 $1085813001,\,1085813004,\,1085813005,\,1085813006,\,1085813007,\,1085813008$ 

Parameter		QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chromatograph	y/Mass S	pectros	сору					
Benzene	LCS	27.9	93	(80-120)			30 ug/L	10/27/2008
Toluene	LCS	29.5	98	(77-120)			30 ug/L	10/27/2008
Ethylbenzene	LCS	30.1	100	(80-120)			30 ug/L	10/27/2008
n-Butylbenzene	LCS	30.2	101	(80-124)			30 ug/L	10/27/2008
1,4-Dichlorobenzene	LCS	29.5	98	(80-120)			30 ug/L	10/27/2008
1,2-Dichloroethane	LCS	28.4	95	(80-129)			30 ug/L	10/27/2008
1,3,5-Trimethylbenzene	LCS	32.7	109	(80-128)			30 ug/L	10/27/2008
4-Chlorotoluene	LCS	31.3	104	(79-128)			30 ug/L	10/27/2008
Chlorobenzene	LCS	30.1	100	(80-120)			30 ug/L	10/27/2008
4-Methyl-2-pentanone (MIBK)	LCS	83.7	93	( 69-134 )			90 ug/L	10/27/2008
cis-1,2-Dichloroethene	LCS	29.3	98	(80-125)			30 ug/L	10/27/2008
4-Isopropyltoluene	LCS	31.8	106	(80-125)			30 ug/L	10/27/2008
cis-1,3-Dichloropropene	LCS	30.2	101	(80-120)			30 ug/L	10/27/2008
n-Propylbenzene	LCS	31.6	105	(80-129)			30 ug/L	10/27/2008
Styrene	LCS	31.7	106	(80-120)			30 ug/L	10/27/2008
Dibromomethane	LCS	29.4	98	(80-120)			30 ug/L	10/27/2008
trans-1,3-Dichloropropene	LCS	32.1	107	(80-124)			30 ug/L	10/27/2008
1,2,4-Trichlorobenzene	LCS	29.0	97	(80-120)			30 ug/L	10/27/2008
Acetone	LCS	81.6	91	(50-135)			90 ug/L	10/27/2008
1,1,2,2-Tetrachloroethane	LCS	30.3	101	(76-123)			30 ug/L	10/27/2008



867630

Lab Control Sample

Printed Date/Time Prep Batch

11/20/2008

8:09

Client Name

The Environmental Company, Inc. (TEC) 9121-003 Red Hill BFSF

Method

VXX18954 SW5030B

Project Name/#

Date 10/27/2008

Matrix	Water	(Surface,	Eff.,	Ground
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Parameter		QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chromatography/M	Mass S	pectrosco	ору					
1,2-Dibromo-3-chloropropane	LCS	34.1	114	(73-130)			30 ug/L	10/27/2008
Methyl-t-butyl ether	LCS	45.0	100	(80-120)			45 ug/L	10/27/2008
Tetrachloroethene	LCS	29.6	99	(79-122)			30 ug/L	10/27/2008
Dibromochloromethane	LCS	31.8	106	(80-120)			30 ug/L	10/27/2008
1,3-Dichloropropane	LCS	30.7	102	(80-121)			30 ug/L	10/27/2008
1,2-Dibromoethane	LCS	31.3	104	(80-120)			30 ug/L	10/27/2008
Carbon tetrachloride	LCS	30.7	102	(80-126)			30 ug/L	10/27/2008
1,1,1,2-Tetrachloroethane	LCS	31.9	106	(80-120)			30 ug/L	10/27/2008
Chloroform	LCS	29.8	99	(80-124)			30 ug/L	10/27/2008
Bromobenzene	LCS	30.8	103	(80-120)			30 ug/L	10/27/2008
Chloromethane	LCS	28.6	95	( 67-125 )			30 ug/L	10/27/2008
1,2,3-Trichloropropane	LCS	30.2	101	(80-120)			30 ug/L	10/27/2008
Bromomethane	LCS	31.6	105	(30-140)			30 ug/L	10/27/2008
Bromochloromethane	LCS	30.5	102	(77-129)			30 ug/L	10/27/2008
Vinyl chloride	LCS	28.8	96	(72-145)			30 ug/L	10/27/2008
Dichlorodifluoromethane	LCS	27.5	92	( 62-153 )			30 ug/L	10/27/2008
Chloroethane	LCS	28.4	95	( 67-133 )			30 ug/L	10/27/2008
sec-Butylbenzene	LCS	32.3	108	(80-120)			30 ug/L	10/27/2008
Bromodichloromethane	LCS	31.8	106	(80-120)			30 ug/L	10/27/2008
1,1-Dichloroethene	LCS	29.5	98	(76-130)			30 ug/L	10/27/2008



867630

Lab Control Sample

Printed Date/Time
Prep Batch

11/20/2008

8:09

Client Name

The Environmental Company, Inc. (TEC)

Batch Method Date VXX18954 SW5030B 10/27/2008

Project Name/#
Matrix

9121-003 Red Hill BFSF

Water (Surface, Eff., Ground)

Parameter		QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chromatography/	Mass S	pectros	сору					
2-Butanone (MEK)	LCS	84.4	94	(66-136)			90 ug/L	10/27/2008
Methylene chloride	LCS	30.2	101	( 63-131 )			30 ug/L	10/27/2008
Trichlorofluoromethane	LCS	30.0	100	( 68-145 )			30 ug/L	10/27/2008
P & M -Xylene	LCS	62.4	104	(80-120)			60 ug/L	10/27/2008
Naphthalene	LCS	28.5	95	(75-120)			30 ug/L	10/27/2008
o-Xylene	LCS	31.5	105	(80-120)			30 ug/L	10/27/2008
Bromoform	LCS	31.8	106	(80-120)			30 ug/L	10/27/2008
1-Chlorohexane	LCS	44.5	99	(70-125)			45 ug/L	10/27/2008
1,2,4-Trimethylbenzene	LCS	31.8	106	(80-125)			30 ug/L	10/27/2008
tert-Butylbenzene	LCS	32.4	108	(80-122)			30 ug/L	10/27/2008
1,1,1-Trichloroethane	LCS	32.0	107	(80-122)			30 ug/L	10/27/2008
1,1-Dichloroethane	LCS	30.6	102	(80-120)			30 ug/L	10/27/2008
2-Chlorotoluene	LCS	31.0	103	(80-125)			30 ug/L	10/27/2008
Trichloroethene	LCS	28.4	95	(80-125)			30 ug/L	10/27/2008
trans-1,2-Dichloroethene	LCS	30.2	101	(79-132)			30 ug/L	10/27/2008
1,2-Dichlorobenzene	LCS	30.4	101	(80-120)			30 ug/L	10/27/2008
2,2-Dichloropropane	LCS	31.2	104	(80-132)			30 ug/L	10/27/2008
Hexachlorobutadiene	LCS	26.4	88	(77-125)			30 ug/L	10/27/2008
Isopropylbenzene (Cumene)	LCS	31.7	106	(80-121)			30 ug/L	10/27/2008
1,2-Dichloropropane	LCS	29.1	97	(80-121)			30 ug/L	10/27/2008
1,1-Dichloropropene	LCS	29.3	98	(80-122)			30 ug/L	10/27/2008



867630

Lab Control Sample

Printed Date/Time

11/20/2008

8:09

Analysis

Date

Prep Batch

Method

RPD

Limits

VXX18954 SW5030B

Spiked

Amount

Date

10/27/2008

Client Name Project Name/# The Environmental Company, Inc. (TEC) 9121-003 Red Hill BFSF

Matrix

Water (Surface, Eff., Ground)

Parameter	Results	Recov	Limits	RPD	
	QC	PCl	LCS/LCSD		

QC

Pct

Volatile Gas Chromatography/Mass Spectroscopy										
1,1,2-Trichloroethane	LCS	30.1	100	(77-120)	30 ug/L	10/27/2008				
1,3-Dichlorobenzene	LCS	29.1	97	(80-120)	30 ug/L	10/27/2008				
1,2,3-Trichlorobenzene	LCS	26.3	88	(77-120)	30 ug/L	10/27/2008				
Surrogates										
1,2-Dichloroethane-D4 <surr></surr>	LCS		103	(73-120)		10/27/2008				
Toluene-d8 <surr></surr>	LCS		104	(80-120)		10/27/2008				
4-Bromofluorobenzene <surr></surr>	LCS		103	(76-120)		10/27/2008				

LCS/LCSD

Batch VMS10242 Method SW8260B

Instrument HP 5890 Series II MS3 VNA



SGS Ref.# 867656 Lab Control Sample Printed Date/Time 11/20/2008 8:09

**Project Name/#** 9121-003 Red Hill BFSF **Date** 10/28/2008

Matrix Water (Surface, Eff., Ground)

QC results affect the following production samples:

 $1085813001,\,1085813004,\,1085813005,\,1085813006,\,1085813007$ 

	QC	Pct	LCS/LCSD		RPD	Spiked	Analysis	
Parameter	Results	Recov	Limits	RPD	Limits	Amount	Date	

Polynuclear Aromatics GC/MS



867656

Lab Control Sample

Printed Date/Time Prep Batch

11/20/2008

8:09

Client Name

The Environmental Company, Inc. (TEC) 9121-003 Red Hill BFSF

Method

XXX20276 SW3520C

Project Name/#

Date 10/28/2008

Matrix	Water (Surface, Eff.,	Ground)
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Parameter		QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Polynuclear Aromatics GC/MS								
Acenaphthylene	LCS	0.333	67	(50-105)			0.5 ug/L	10/30/2008
Acenaphthene	LCS	0.314	63	(45-110)			0.5 ug/L	10/30/2008
Fluorene	LCS	0.342	68	(50-110)			0.5 ug/L	10/30/2008
Phenanthrene	LCS	0.369	74	(50-115)			0.5 ug/L	10/30/2008
Anthracene	LCS	0.382	76	(55-110)			0.5 ug/L	10/30/2008
Fluoranthene	LCS	0.342	68	(55-125)			0.5 ug/L	10/30/2008
Pyrene	LCS	0.323	65	(50-130)			0.5 ug/L	10/30/2008
Benzo(a)Anthracene	LCS	0.376	75	(55-120)			0.5 ug/L	10/30/2008
Chrysene	LCS	0.339	68	(55-120)			0.5 ug/L	10/30/2008
Benzo[b]Fluoranthene	LCS	0.376	75	( 46-130 )			0.5 ug/L	10/30/2008
Benzo[k]fluoranthene	LCS	0.378	76	( 60-125 )			0.5 ug/L	10/30/2008
Benzo[a]pyrene	LCS	0.416	83	(55-120)			0.5 ug/L	10/30/2008
Indeno[1,2,3-c,d] pyrene	LCS	0.379	76	(45-125)			0.5 ug/L	10/30/2008
Dibenzo[a,h]anthracene	LCS	0.389	78	(41-140)			0.5 ug/L	10/30/2008
Benzo[g,h,i]perylene	LCS	0.376	75	(46-125)			0.5 ug/L	10/30/2008
Naphthalene	LCS	0.266	53	( 42-100 )			0.5 ug/L	10/30/2008
1-Methylnaphthalene	LCS	0.277	55	(46-115)			0.5 ug/L	10/30/2008
2-Methylnaphthalene	LCS	0.290	58	( 45-105 )			0.5 ug/L	10/30/2008
Surrogates								
Terphenyl-d14 <surr></surr>	LCS		93	(50-135)				10/30/2008



 SGS Ref.#
 867656
 Lab Control Sample
 Printed Date/Time
 11/20/2008
 8:09

 Prep
 Batch
 XXX20276

Client Name The Environmental Company, Inc. (TEC) Method SW3520C

Project Name/#9121-003 Red Hill BFSFDate10/28/2008MatrixWater (Surface, Eff., Ground)

Parameter QC Pct LCS/LCSD RPD Spiked Analysis
Recov Limits RPD Limits Amount Date

## Polynuclear Aromatics GC/MS

BatchXMS4761Method8270D SIMS

Instrument HP 5890 Series II MS2 SVOA



Client Name

SGS Ref.# 868152 Lab Control Sample

868153 Lab Control Sample Duplicate

The Environmental Company, Inc. (TEC)

Project Name/# 9121-003 Red Hill BFSF

Matrix Water (Surface, Eff., Ground)

Printed Date/Time

Prep

11/20/2008

Batch Method XXX20290 SW3520C 8:09

Date 10/29/2008

QC results affect the following production samples:

1085813004, 1085813005, 1085813006, 1085813007

Parameter	QC Resul		LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Semivolatile Organic Fu	els Department						
Diesel Range Organics	LCS 4.42	89	(75-125)			5 mg/L	10/31/2008
	LCSD 5.32	106	,	18	(< 20)	5 mg/L	10/31/2008
Surrogates							
5a Androstane <surr></surr>	LCS	100	(60-120)				10/31/2008
	LCSD	110		10			10/31/2008

Batch XFC8323 Method SW8015C

Instrument HP 6890 Series II FID SV D R



Client Name

Matrix

868325 Lab Control Sample

868326 Lab Control Sample Duplicate

The Environmental Company, Inc. (TEC)

Project Name/#

9121-003 Red Hill BFSF

Water (Surface, Eff., Ground)

Printed Date/Time

Prep

Batch VXX18967

Method

SW5030B

11/20/2008

8:09

Date

10/30/2008

QC results affect the following production samples:

 $1085813001,\,1085813004,\,1085813005,\,1085813006,\,1085813007$ 

Parameter		QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Fuels Department								
Gasoline Range Organics	LCS	192	96	(79-108)			200 ug/L	10/30/2008
	LCSD	189	95		2	(< 20)	200 ug/L	10/30/2008
Surrogates								
4-Bromofluorobenzene <surr></surr>	LCS		96	(50-150)				10/30/2008
	LCSD		95		0			10/30/2008

Batch VFC9248 Method SW8015C

Instrument HP 5890 Series II PID+FID VCA



Matrix

868747 Lab Control Sample

868748 Lab Control Sample Duplicate

The Environmental Company, Inc. (TEC)

Client Name Project Name/#

9121-003 Red Hill BFSF

Water (Surface, Eff., Ground)

Printed Date/Time

Prep

11/20/2008

8:09

Batch Method VXX18976 SW5030B

Date

10/31/2008

QC results affect the following production samples:

 $1085813005,\,1085813006,\,1085813007$ 

Parameter	QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chromatography,	Mass Spectros	сору					
Naphthalene	LCS 29.2	97	(75-120)			30 ug/L	10/31/2008
	LCSD 28.9	96		1	(< 20)	30 ug/L	10/31/2008
Surrogates							
1,2-Dichloroethane-D4 <surr></surr>	LCS	102	(73-120)				10/31/2008
	LCSD	99		3			10/31/2008
Toluene-d8 <surr></surr>	LCS	99	(80-120)				10/31/2008
	LCSD	99		0			10/31/2008
4-Bromofluorobenzene <surr></surr>	LCS	99	(76-120)				10/31/2008
	LCSD	97		2			10/31/2008

Batch VMS10254 Method SW8260B

Instrument HP 5890 Series II MS3 VNA



Client Name

Matrix

Project Name/#

869198

Lab Control Sample

The Environmental Company, Inc. (TEC)

Printed Date/Time

11/20/2008

8:09

Prep Batch

XXX20313

Method Date SW3520C 11/03/2008

9121-003 Red Hill BFSF Water (Surface, Eff., Ground)

QC results affect the following production samples:

1085813001

Parameter	QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Semivolatile Organic Fuels	Department						
Diesel Range Organics	LCS 2.81	56 *	(75-125)			5 mg/L	11/05/2008
Surrogates							
5a Androstane <surr></surr>	LCS	84	(60-120)				11/05/2008

Batch XFC8338 Method SW8015C

**Instrument** HP 6890 Series II FID SV D R



SGS Ref.# 869582 Lab Control Sample Printed Date/Time

The Environmental Company, Inc. (TEC)

Printed Date/Time 11/20/2008 Prep Batch MXX21031

Method SW3010A

8:09

9121-003 Red Hill BFSF **Date** 11/04/2008

Matrix Water (Surface, Eff., Ground)

QC results affect the following production samples:

 $1085813001,\, 1085813004,\, 1085813005,\, 1085813006$ 

Parameter QC Pct LCS/LCSD RPD Spiked Analysis
Results Recov Limits RPD Limits Amount Date

Metals by ICP/MS

Client Name

Project Name/#

Lead LCS 1070 107 (80-120) 1000 ug/L 11/14/2008

Batch MMS5719 Method SW6020

**Instrument** Perkin Elmer Sciex ICP-MS P4



868545 868546 Matrix Spike

Matrix Spike Duplicate

Printed Date/Time

11/20/2008 8:09

Prep Batch XXX20290

Method

Cnt. Liq/Liq Ext. for AK102/3

Date 10/29/2008

Original 871926

Matrix

Water (Surface, Eff., Ground)

QC results affect the following production samples:

 $1085813004,\,1085813005,\,1085813006,\,1085813007$ 

Parameter	Qualifiers	Original Result	QC Result	Pct Recov	MS/MSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Semivolatile Org	anic Fuels	Departmen	<u>t</u>						
Diesel Range Organics	MS MSI	ND	ND 0.00	0* 0*	(75-125)	0	(< 30)	0.20	g/L 10/31/2008 g/L 10/31/2008
Surrogates	11101		0.00	v		v	( 50)	3.20 m	.g/ E 10/31/2000
5a Androstane <surr></surr>	MS		.107	102	(50-150)				10/31/2008
	MSI	)	0.102	97		5			10/31/2008
Batch XF0	C8323								

Method SW8015C

Instrument HP 6890 Series II FID SV D R



 $1085813002 \\ 1085813003$ 

Billable Matrix Spike

Billable Matrix Spike Dup.

Printed Date/Time

11/20/2008 8:09

Prep Batch MXX21031

Date

Method 3010 H20 Digest for Metals ICI

11/04/2008

Original

1085813001

Matrix Water (Surface, Eff., Ground)

QC results affect the following production samples:

Parameter	Qualifier	Original rs Result	QC Result	Pct Recov	MS/MSD Limits	RPD	RPD Limits	Spiked Amount	
Dissolved M	Metals by ICP	/MS							
Lead		BMS ND	1020	102	(80-120)			1000	ug/L 11/14/2008
		BMSD	1000	100		1	(< 15)	1000	ug/L 11/14/2008
Batch	MMS5719								
Method	SW6020								
Instrument	Perkin Elmer	Sciex ICP-MS P4							
/olatile Fu	els Departme	<u>nt</u>							
Gasoline Range	Organics	BMS ND	448	100	(79-108)			450	ug/L 10/30/2008
		BMSD	401	89		11	(< 20)	450	ug/L 10/30/2008
Surrogates									
1-Bromofluorob	enzene <surr></surr>	BMS	50.9	102	(50-150)				10/30/2008
		BMSD	48.6	97		5			10/30/2008
Batch	VFC9248								
Method	SW8015C								
Instrument	HP 5890 Seri	es II PID+FID VCA	L						
Semivolatil	e Organic Fu	els Department	<u> </u>						
Diesel Range Or	rganics	BMS ND	4.89	94	(75-125)			5.21	mg/L 11/05/2008
		BMSD	4.72	94		4	(< 30)	5.00	mg/L 11/05/2008
Surrogates									
5a Androstane <	<surr></surr>	BMS	.103	99	(50-150)				11/05/2008
		BMSD	0.102	102		1			11/05/2008
Batch Method	XFC8338 SW8015C								

# Volatile Gas Chromatography/Mass Spectroscopy

HP 6890 Series II FID SV D R

Instrument



 $\frac{1085813002}{1085813003}$ 

Billable Matrix Spike

Billable Matrix Spike Dup.

Printed Date/Time
Prep Batch

ne 11/20/2008 8:09 ch VXX18954

Batch VXX18954
Method Volatiles Extraction AFCEE 3.1

Date 10/27/2008

**Original** 1085813001

Parameter	Qualifiers	Original Result	QC Result	Pct Recov	MS/MSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chr	omatography,	/Mass Spe	ctroscopy						
Benzene	BMS	S ND	27.9	93	(80-120)			30.0	ug/L 10/27/2008
	BMS	SD	28.4	95		2	(< 20)		ug/L 10/27/2008
Toluene	BMS	S ND	29.6	99	(77-120)				ug/L 10/27/2008
	BMS	SD	29.6	99		0	(< 20)		ug/L 10/27/2008
Ethylbenzene	BMS	S ND	30.6	102	(80-120)				ug/L 10/27/2008
•	BMS	SD	31.3	104		2	(< 20)		ug/L 10/27/2008
n-Butylbenzene	BMS	S ND	34.4	115	(80-124)				ug/L 10/27/2008
•	BMS		35.5	118		3	(< 20)		ug/L 10/27/2008
1,4-Dichlorobenzene	BMS	S ND	29.3	98	(80-120)				ug/L 10/27/2008
•	BMS		30.9	103		5	(< 20)		ug/L 10/27/2008
1,2-Dichloroethane	BMS	S ND	28.3	94	(80-129)				ug/L 10/27/2008
	BMS		29.2	97		3	(< 20)		ug/L 10/27/2008
1,3,5-Trimethylbenzen		S ND	33	110	(80-128)		,		ug/L 10/27/2008
•	BMS		34.0	113		3	(< 20)		ug/L 10/27/2008
4-Chlorotoluene	BMS	S ND	31.8	106	(79-128)				ug/L 10/27/2008
	BMS		33.0	110		4	(< 20)		ug/L 10/27/2008
Chlorobenzene		S ND	30	100	(80-120)		,		ug/L 10/27/2008
	BMS		29.9	100	, ,	0	(< 20)		ug/L 10/27/2008
4-Methyl-2-pentanone		S ND	82.4	92	(69-134)		,		ug/L 10/27/2008
J P	BMS		87.9	98	,	7	(< 20)		ug/L 10/27/2008
cis-1,2-Dichloroethene		S ND	29.4	98	(80-125)		,		ug/L 10/27/2008
,	BMS		29.8	99	,	1	(< 20)		ug/L 10/27/2008
4-Isopropyltoluene		S ND	33.7	112	(80-125)		,		ug/L 10/27/2008
1 13	BMS		34.4	115	, , ,	2	(< 20)		ug/L 10/27/2008
cis-1,3-Dichloroproper		S ND	29.5	98	(80-120)		,		ug/L 10/27/2008
, , , , , , , , , , , , , , , , , , ,	BMS		30.8	103	,	5	(< 20)		ug/L 10/27/2008
n-Propylbenzene		S ND	32.5	108	(80-129)		,		ug/L 10/27/2008
13	BMS		33.7	112	, , ,	4	(< 20)		ug/L 10/27/2008
Styrene		S ND	32	107	(80-120)		,		ug/L 10/27/2008
	BMS		32.3	108	,	1	(< 20)		ug/L 10/27/2008
Dibromomethane		S ND	28.4	95	(80-120)		,		ug/L 10/27/2008
	BMS		29.7	99	, ,	5	(< 20)		ug/L 10/27/2008
trans-1,3-Dichloroprop		S ND	30.1	100	(80-124)		,		ug/L 10/27/2008
, , , , , , , , , , , , , , , , , , ,	BMS		31.5	105	,	5	(< 20)		ug/L 10/27/2008
1,2,4-Trichlorobenzene		S ND	30.9	103	(80-120)		, ,		ug/L 10/27/2008
, ,	BMS		32.2	107	. ,	4	(< 20)		ug/L 10/27/2008
Acetone		S ND	83.2	93	(50-135)		, ,		ug/L 10/27/2008
	BMS		86.9	97	,	4	(< 20)		ug/L 10/27/2008
1,1,2,2-Tetrachloroetha		S ND	29.3	98	(76-123)		` /		ug/L 10/27/2008
, ,=,=	BMS		30.7	102	- /	4	(< 20)		ug/L 10/27/2008



 $\frac{1085813002}{1085813003}$ 

Billable Matrix Spike

Billable Matrix Spike Dup.

Printed Date/Time
Prep Batch

e/Time 11/20/2008 8:09 Batch VXX18954

Method Volatiles Extraction AFCEE 3.1

Date 10/27/2008

**Original** 1085813001

Volatile Gas Chronological 1,2-Dibromo-3-chloroprological Methyl-t-butyl ether  Tetrachloroethene  Dibromochloromethane  1,3-Dichloropropane  1,2-Dibromoethane	pane Bi Bi Bi Bi Bi Bi Bi	MS ND MSD	32.6 34.4 45.4 47.2 30.6 30.5 29.9	109 115 101 105 102 102	(73-130) (80-120) (79-122)	5	(< 20 )	30.0 45.0	ug/L 10/27/2008 ug/L 10/27/2008 ug/L 10/27/2008
Methyl-t-butyl ether Tetrachloroethene Dibromochloromethane 1,3-Dichloropropane	Bi Bi Bi Bi Bi Bi	MSD MS ND MSD MS ND MSD MS ND MSD MS ND MSD	34.4 45.4 47.2 30.6 30.5	115 101 105 102	(80-120)			30.0 45.0	ug/L 10/27/2008
Methyl-t-butyl ether Tetrachloroethene Dibromochloromethane 1,3-Dichloropropane	Bi Bi Bi Bi Bi Bi	MS ND MSD MS ND MSD MS ND MSD MS ND MSD	45.4 47.2 30.6 30.5	101 105 102				30.0 45.0	ug/L 10/27/2008
Tetrachloroethene Dibromochloromethane 1,3-Dichloropropane	B! B! B! B! B!	MSD MS ND MSD MS ND MSD	45.4 47.2 30.6 30.5	105 102		4		45.0	•
Tetrachloroethene Dibromochloromethane 1,3-Dichloropropane	BM BM BM BM BM	MS ND MSD MS ND MSD	30.6 30.5	102	(79-122)	4	(< 20.)		
Dibromochloromethane 1,3-Dichloropropane	Bi Bi Bi	MSD MS ND MSD	30.5		(79-122)		(< 20)	45.0	ug/L 10/27/2008
Dibromochloromethane 1,3-Dichloropropane	Bi Bi Bi	MSD MS ND MSD	30.5		(1) 144 )				ug/L 10/27/2008
1,3-Dichloropropane	Bi Bi	MSD				0	(< 20)		ug/L 10/27/2008
	BI			100	(80-120)				ug/L 10/27/2008
		AC NE	30.7	102		3	(< 20)		ug/L 10/27/2008
		MS ND	30.2	101	(80-121)				ug/L 10/27/2008
1,2-Dibromoethane		MSD	30.6	102		1	(< 20)		ug/L 10/27/2008
	BN	MS ND	29.8	99	(80-120)				ug/L 10/27/2008
	Bl	MSD	30.4	101		2	(< 20)		ug/L 10/27/2008
Carbon tetrachloride		MS ND	29.4	98	(80-126)		,		ug/L 10/27/2008
		MSD	30.3	101		3	(< 20)		ug/L 10/27/2008
1,1,1,2-Tetrachloroethan	BI	MS ND	30.2	101	(80-120)				ug/L 10/27/2008
, , ,		MSD	30.6	102		1	(< 20)		ug/L 10/27/2008
Chloroform		MS ND	29.2	97	(80-124)		,		ug/L 10/27/2008
		MSD	29.7	99		2	(< 20)		ug/L 10/27/2008
Bromobenzene		MS ND	29.9	100	(80-120)		,		ug/L 10/27/2008
		MSD	31.2	104	,	4	(< 20)		ug/L 10/27/2008
Chloromethane		MS ND	28.2	94	(67-125)		,		ug/L 10/27/2008
		MSD	28.9	96	,	2	(< 20)		ug/L 10/27/2008
1,2,3-Trichloropropane		MS ND	30.4	101	(80-120)		,		ug/L 10/27/2008
, ,- · · · · · · · · · · · · · · · · · ·		MSD	31.4	105	,	3	(< 20)		ug/L 10/27/2008
Bromomethane		MS ND	32.8	109	(30-140)		,		ug/L 10/27/2008
		MSD	35.7	119	,	9	(< 20)		ug/L 10/27/2008
Bromochloromethane		MS ND	28.4	95	(77-129)		,		ug/L 10/27/2008
		MSD	29.0	97	,	2	(< 20)		ug/L 10/27/2008
Vinyl chloride		MS ND	29.6	99	(72-145)				ug/L 10/27/2008
,, . <b>.</b>		MSD	30.3	101	( ' ' ' '	2	(< 20)		ug/L 10/27/2008
Dichlorodifluoromethane		MS ND	28.7	96	(62-153)				ug/L 10/27/2008
		MSD	29.1	97	,	2	(< 20)		ug/L 10/27/2008
Chloroethane		MS ND	30	100	(67-133)	_	( = ,		ug/L 10/27/2008
		MSD	34.5	115	,	14	(< 20)		ug/L 10/27/2008
sec-Butylbenzene		MS ND	33.3	111	(80-120)		)		ug/L 10/27/2008
		MSD	34.7	116		4	(< 20)		ug/L 10/27/2008
Bromodichloromethane		MS ND	30.5	102	(80-120)	-	· · /		ug/L 10/27/2008
		MSD	31.8	106	/	4	(< 20)		ug/L 10/27/2008
1,1-Dichloroethene		MS ND	31.6	105	(76-130)	·			ug/L 10/27/2008
1,1 Diemoroemene		MSD	32.5	108	( )	3	(< 20)		ug/L 10/27/2008



1085813002 1085813003 Billable Matrix Spike

Billable Matrix Spike Dup.

Printed Date/Time Prep Batch

11/20/2008 8:09 VXX18954

Batch VXX18954 Method Volatiles E

**Date** 10/2

Volatiles Extraction AFCEE 3.1 10/27/2008

**Original** 1085813001

Parameter		ginal QC sult Result	Pct Recov	MS/MSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chrom	natography/Mass	Spectroscopy						
2-Butanone (MEK)	BMS ND	85.1	95	(66-136)			90.0	ug/L 10/27/2008
, ,	BMSD	87.2	97		2	(< 20)		ug/L 10/27/2008
Methylene chloride	BMS ND		104	(63-131)		,		ug/L 10/27/2008
,	BMSD	32.9	110	,	5	(< 20)		ug/L 10/27/2008
Trichlorofluoromethane	BMS ND		102	(68-145)		,		ug/L 10/27/2008
	BMSD	31.0	103		1	(< 20)		ug/L 10/27/2008
P & M -Xylene	BMS ND		108	(80-120)				ug/L 10/27/2008
<b>y</b>	BMSD	65.1	109	,	1	(< 20)		ug/L 10/27/2008
Naphthalene	BMS ND		96	(75-120)		,		ug/L 10/27/2008
	BMSD	30.2	101	,	5	(< 20)		ug/L 10/27/2008
o-Xylene	BMS ND		106	(80-120)		,		ug/L 10/27/2008
. ,	BMSD	31.8	106	,	0	(< 20)		ug/L 10/27/2008
Bromoform	BMS ND		100	(80-120)				ug/L 10/27/2008
	BMSD	30.9	103	,	3	(< 20)		ug/L 10/27/2008
1-Chlorohexane	BMS ND		111	(70-125)				ug/L 10/27/2008
T Children Children	BMSD	50.2	112	( )	0	(< 20)		ug/L 10/27/2008
1,2,4-Trimethylbenzene	BMS ND		108	(80-125)	-	( = - )		ug/L 10/27/2008
1,2,1 1111110111111001120110	BMSD	33.6	112	( ** -=* )	4	(< 20)		ug/L 10/27/2008
tert-Butylbenzene	BMS ND		107	(80-122)		( = - )		ug/L 10/27/2008
tert Butyroenzene	BMSD	33.6	112	(00 122)	5	(< 20)		ug/L 10/27/2008
1,1,1-Trichloroethane	BMS ND		104	(80-122)	J	( '20 )		ug/L 10/27/2008
1,1,1 Themoreemane	BMSD	32.1	107	(00 122)	3	(< 20)		ug/L 10/27/2008
1,1-Dichloroethane	BMS NE		105	(80-120)	3	( 120 )		ug/L 10/27/2008
1,1 Diemorochane	BMSD	32.4	108	(00120)	3	(< 20)		ug/L 10/27/2008
2-Chlorotoluene	BMS ND		103	(80-125)	J	( '20 )		ug/L 10/27/2008
2 Chiorototache	BMSD	32.2	107	(00 120)	4	(< 20)		ug/L 10/27/2008
Trichloroethene	BMS ND		97	(80-125)	•	( '20 )		ug/L 10/27/2008
THEMOTOCHIC	BMSD	29.3	98	(00 120)	1	(< 20)		ug/L 10/27/2008
trans-1,2-Dichloroethene	BMS ND		109	(79-132)	•	( '20 )		ug/L 10/27/2008
trans-1,2-Diemorocurene	BMSD	33.0	110	(7) 132)	1	(< 20)		ug/L 10/27/2008 ug/L 10/27/2008
1,2-Dichlorobenzene	BMS ND		101	(80-120)	1	( 120 )		ug/L 10/27/2008
1,2-Diemoroochzene	BMSD	31.4	105	(00 120)	4	(< 20)		ug/L 10/27/2008 ug/L 10/27/2008
2,2-Dichloropropane	BMS ND		106	(80-132)	7	( \ 20 )		ug/L 10/27/2008 ug/L 10/27/2008
2,2-Diemoropropane	BMSD	32.9	110	(00 132)	3	(< 20)		ug/L 10/27/2008 ug/L 10/27/2008
Hexachlorobutadiene	BMS NE		99	(77-125)	3	( 120 )		ug/L 10/27/2008 ug/L 10/27/2008
ricacinorodulautene	BMSD	30.0	100	( / / 123 )	1	(< 20)		ug/L 10/27/2008 ug/L 10/27/2008
Isopropylbenzene (Cumer			100	(80-121)	1	( 20 )		ug/L 10/27/2008 ug/L 10/27/2008
150propyrochizene (Cumer	BMSD		108	(00-121)	1	(< 20)		ug/L 10/27/2008 ug/L 10/27/2008
1,2-Dichloropropane		32.7 28.9	96	(80-121)	1	( > 20 )		-
1,2-Dichioropropane	BMS ND			(00-121)	2	(< 20 )		ug/L 10/27/2008
	BMSD	29.8	99		3	(< 20)	30.0	ug/L 10/27/2008



1085813002 1085813003 Billable Matrix Spike

Billable Matrix Spike Dup.

Printed Date/Time

Prep

11/20/2008 8:09

VXX18954

Batch Method Volatiles Extraction AFCEE 3.1

Date 10/27/2008

Original

1085813001

Matrix Water (Surface, Eff., Ground)

		.,	)						
Parameter (	Qualifiers	Original Result	QC Result	Pct Recov	MS/MSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chrom	atography/	Mass Spe	ctroscopy						
1,1-Dichloropropene	BMS	ND	30.3	101	(80-122)			30.0 u	ıg/L 10/27/2008
	BMS	D	30.5	102		1	(< 20)	30.0 ι	ıg/L 10/27/2008
1,1,2-Trichloroethane	BMS	ND	29.7	99	(77-120)			30.0 u	ıg/L 10/27/2008
	BMS	D	30.1	100		2	(< 20)	30.0 ι	ıg/L 10/27/2008
1,3-Dichlorobenzene	BMS	ND	29.5	99	(80-120)			30.0 u	ıg/L 10/27/2008
	BMS	D	30.5	102		3	(< 20)	30.0 ι	ıg/L 10/27/2008
1,2,3-Trichlorobenzene	BMS	ND	27.7	92	(77-120)			30.0 u	ıg/L 10/27/2008
	BMS	D	29.3	98		5	(< 20)	30.0 u	ıg/L 10/27/2008
Surrogates									
1,2-Dichloroethane-D4 <s< td=""><td>urr&gt; BMS</td><td></td><td>30.9</td><td>103</td><td>(73-120)</td><td></td><td></td><td></td><td>10/27/2008</td></s<>	urr> BMS		30.9	103	(73-120)				10/27/2008
	BMS	D	31.0	103		0			10/27/2008
Toluene-d8 <surr></surr>	BMS		30.9	103	(80-120)				10/27/2008
	BMS	D	30.8	103		0			10/27/2008
4-Bromofluorobenzene <s< td=""><td>urr&gt; BMS</td><td></td><td>30.3</td><td>101</td><td>(76-120)</td><td></td><td></td><td></td><td>10/27/2008</td></s<>	urr> BMS		30.3	101	(76-120)				10/27/2008
	BMS	D	31.0	103		2			10/27/2008

Batch VMS10242 Method SW8260B

Instrument HP 5890 Series II MS3 VNA

Polynuclear Aromatics GC/MS



 $\frac{1085813002}{1085813003}$ 

Billable Matrix Spike

Billable Matrix Spike Dup.

Printed Date/Time
Prep Batch

11/20/2008 8:09 XXX20276

**Batch** XXX20276 **Method** 3520 Liquid/Liquid Ext for 827

Date 10/28/2008

**Original** 1085813001

Parameter	Qualifiers	Original Result	QC Result	Pct Recov	MS/MSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Del	-ti 00/20								
Polynuclear Arom	atics GC/MS								
Acenaphthylene		S ND	.362	68	(50-105)			0.532	ug/L 10/30/2008
	BMS		0.350	64		3	(< 30)	0.549	ug/L 10/30/2008
Acenaphthene		S ND	.345	65	(45-110)			0.532	ug/L 10/30/2008
	BMS	SD	0.342	62		1	(< 30)	0.549	ug/L 10/30/2008
Fluorene		S ND	.36	68	(50-110)			0.532	ug/L 10/30/2008
	BMS		0.349	64		3	(< 30)	0.549	ug/L 10/30/2008
Phenanthrene		S ND	.348	65	(50-115)			0.532	ug/L 10/30/2008
	BMS	SD	0.343	63		1	(< 30)	0.549	ug/L 10/30/2008
Anthracene		S ND	.372	70	(55-110)			0.532	ug/L 10/30/2008
	BMS	SD	0.363	66		2	(< 30)	0.549	ug/L 10/30/2008
Fluoranthene		S ND	.38	71	(55-125)			0.532	ug/L 10/30/2008
	BMS	SD	0.378	69		1	(< 30)	0.549	ug/L 10/30/2008
Pyrene		S ND	.366	69	(50-130)			0.532	ug/L 10/30/2008
	BMS	SD	0.364	66		1	(< 30)	0.549	ug/L 10/30/2008
Benzo(a)Anthracene	BMS	S ND	.398	75	(55-120)			0.532	ug/L 10/30/2008
	BMS	SD	0.380	69		5	(< 30)	0.549	ug/L 10/30/2008
Chrysene	BMS	S ND	.386	73	(55-120)			0.532	ug/L 10/30/2008
	BMS	SD	0.387	70		0	(< 30)	0.549	ug/L 10/30/2008
Benzo[b]Fluoranthene	BMS	S ND	.406	76	(46-130)			0.532	ug/L 10/30/2008
	BMS	SD	0.408	74		1	(< 30)	0.549	ug/L 10/30/2008
Benzo[k]fluoranthene	BMS	S ND	.413	78	(60-125)			0.532	ug/L 10/30/2008
	BMS	SD	0.392	71		5	(< 30)	0.549	ug/L 10/30/2008
Benzo[a]pyrene	BMS	S ND	.43	81	(55-120)			0.532	ug/L 10/30/2008
	BMS	SD	0.418	76		3	(< 30)	0.549	ug/L 10/30/2008
Indeno[1,2,3-c,d] pyren	e BMS	S ND	.416	78	(45-125)			0.532	ug/L 10/30/2008
	BMS	SD	0.414	75		1	(< 30)	0.549	ug/L 10/30/2008
Dibenzo[a,h]anthracene	e BMS	S ND	.424	80	(41-140)			0.532	ug/L 10/30/2008
	BMS	SD	0.425	77		0	(< 30)	0.549	ug/L 10/30/2008
Benzo[g,h,i]perylene	BMS	S ND	.421	79	(46-125)			0.532	ug/L 10/30/2008
	BMS	SD	0.416	76		1	(< 30)	0.549	ug/L 10/30/2008
Naphthalene	BMS	S 0.0466 J	.333	54	(42-100)			0.532	ug/L 10/30/2008
	BMS	SD	0.325	51		3	(< 30)	0.549	ug/L 10/30/2008
1-Methylnaphthalene	BMS	S 0.0276 J	.341	59	(46-115)			0.532	ug/L 10/30/2008
	BMS	SD	0.336	56		2	(< 30)	0.549	ug/L 10/30/2008
2-Methylnaphthalene	BMS	S ND	.336	63	(45-105)			0.532	ug/L 10/30/2008
	BMS	SD	0.327	60		3	(< 30)	0.549	ug/L 10/30/2008
Surrogates									
Terphenyl-d14 <surr></surr>	BMS	S	.457	86	(50-135)				10/30/2008
	BMS		0.454	83	. ,	1			10/30/2008



1085813002 1085813003

Billable Matrix Spike

Billable Matrix Spike Dup.

Printed Date/Time

11/20/2008 8:09

Prep Batch XXX20276

Date

Method 3520 Liquid/Liquid Ext for 827

10/28/2008

Original

1085813001

Matrix

Water (Surface, Eff., Ground)

Pct MS/MSD RPD QC Spiked Analysis Original Qualifiers RPD Parameter Limits Recov Limits Amount Result Result Date

## Polynuclear Aromatics GC/MS

Batch XMS4761 Method 8270D SIMS

Instrument HP 5890 Series II MS2 SVOA

# Hager, Barbara (Anchorage)

From: Adkisson, Richard K. [RKAdkisson@tecinc.com]

Sent: Wednesday, October 29, 2008 8:58 AM

To: Hager, Barbara (Anchorage)

Cc: Hart, Jeff; Whitman, William M.C.; Adkisson, Richard K.

**Subject:** Analysis of Sample RHMW01-WG13

Importance: High

Attachments: image001.jpg

# Barbara,

Please <u>cancel</u> the analysis of sample RHMW01-WG13 <u>for total lead</u>. TEC will collect another sample, filter it in the field, and sent it for <u>dissolved</u> lead analysis. Please press on with the other requested analysis for RHMW01-WG13 and all other samples. Thank you.



Rick Adkisson TEC, Inc. 1001 Bishop Street, Suite 1400 Honolulu, HI 96813

Phone: 808-528-1445 Fax: 808-528-0768

# Hager, Barbara (Anchorage)

From:

Adkisson, Richard K. [RKAdkisson@tecinc.com]

Sent:

Thursday, October 23, 2008 9:11 AM

To:

Hager, Barbara (Anchorage)

Cc:

Adkisson, Richard K.; Whitman, William M.C.

Subject:

RE: GW Sampling 10/22 - SPECIAL REQUEST®

Importance:

High

Attachments: image001.jpg

Barbara,

With the Red Hill GW samples that you should be receiving either today or tomorrow you need to be aware of the following:

• Regarding GW Sample RHMW01-WG13 to be analyzed for dispolyed lead: We were <u>not</u> able to filter this particular GW sample in the field prior to shipment, therefore <u>BEFORE</u> analyzing this <u>one</u> sample please filter it in the lab. All other GW samples for dissolved lead <u>have been</u> field filtered.

Also the GW Sample RHMW01-WG13, we were only able to collect 1 (one) vs. 2 (two) amber jars (unpreserved)
for this sample due to low pump flow. Mark Abe verified that your lab <u>only</u> requires 1 (one) jar for its analysis
and that the other jar is requested in the event of breakage.

Please reply by email that you have received this special request ASAP. Also, we <u>cannot afford</u> to have the types of delays and mistakes in the analytical data and it's reporting to us that we recently experienced. Our client has been pressing us to ensure that <u>from now on</u> the analytical data will be presented in a <u>timely and accurate manner</u>.

Barbara, please take personal ownership with this batch of samples to ensure that TEC and our client receive professional, top-notch services.

Thank you and please an knowledge your receipt of this email.



Rick Adkisson TEC, Inc. 1001 Bishop Street, Suite 1400 Honolulu, HI 96813 Phone: 808-528-1445

From: Whitman, William M.C.

**Sent:** Tuesday, October 21, 2008 10:54 AM

Fax: 808-528-0768

**To:** Hager, Barbara (Anchorage) **Cc:** Adkisson, Richard K.; Hart, Jeff **Subject:** GW Sampling 10/22

Hi Barbara,

1085813

10/24/2008

10:::4 200a





#### Locations Nationwide

Alaska

Hawaii

West Virginia

Maryland

Louisiana

New Jersey

North Carolina

WWW.US.SQS.COM

		<u> </u>			SGS Refe	erence #:															7
CLIENT:	TEC INC.				•												page	·	(	of	
CONTACT:	Jeff Hart	PHONE NO:	303,273,0231			Preserv.	,	,				, ,			إ	<u> </u>	_		, ,		-
PROJECT:	9121-003	SITE/PWSID#:	Red Hill BFSF			Used	ķū	$\angle$	, KC	$\angle$	NY.	/	_	_	_	_	_	_			
REPORTS TO:	Jeff Hart	email <u>jshart@tec</u> cc wmcwhitm	<u>inc.com</u> an@tecinc.co	m	°	SAMPLE TYPE C =			-	IS)								1			
INVOICE TO:		QUOTE #:			N T A I	COMP	TPH-GRO (8015B)	TPH-DRO (8015B)	8260B)	(8270C-SIMS)	Pb (6020)		7	دع	3052	86 X	E B	200	240	B	
LAB (40.	SAMPLE IDENTIFICATION	DATE	TIME	MATRIX	E R. 9	G = GRAB	TPH-GR	TPH-DR	VOC's (8260B)	PAH's (	Diss Pb		X				K			REMARKS_	
n Kara	RHMW2254-WG13	10/22/2008	0930	Water	23		Х	X	X	X	X		) (	D/	3)				3x Volum	ne sent in 2 coole	2
4) A - F	RHMW03-WG13	10/22/2008	1115	Water	6		Х		X			) _	٦.	X	6						
5)	RHMW02-WG13	10/22/2008	1245	Water	6		Х		X				<u> </u>	4	χ,						
9A-F 5) 7 V 5A-C	RHMWA01-WG13	10/22/2008	1205	Water	6		Х		X					No.	*	个					
<b>₹</b>	RHMW01-WG13	10/22/2008	1500	Water	6_		Х		X			18 E.	Ĺ	5	14	'					
\$A-C	TB01-WG13	10/22/2008	0805	Water	3				X	1 (C.)	1	, ¥.	Me.		1						
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Callected/Relinquis	hed By: (1)	Date 10-22-08	1720	Received By:		P	1	フ		ng Carri							1 77	7		YES NO ちしてニシ	6
Relinquished By	(2)	Date	Time	Received By:			-					Requiren	nents:				Chain	of Cu	stody Seal:	(Circle)	
/		10/23/03	1435						See (	Contrac	:t						INTAC	T 8	ROKEN	ABSENT	
Relinquished By: (3)  Date Time Received B											marour n <b>trac</b>		end-o	r Speci	ial Instr	uctions	:	·			
Relinquished By:	(4)	Date 1/24/04	Time 1050	Received For Lab	pratory By:	)										<u></u>			·	· <del>·</del>	
3180 Pege	otter Drive Anchorage, AK 99518 Tel: r Road Fairbanks, AK 99701 Tel: (907 Island Access Rd., Unit 1B Honolulu, H	) 474-8656 Fax: (907) 47-	4-9685				Green	brier St	reet Ch	arlesto	n, WV	25311	Tel: (	304) 34	6401 F 16-0725 1903	Fax	(304)	346-07			



1085813

#### Locations Nationwide

Alaska

Hawaii

Maryland

Louisiana

1017 301309

New Jersey West Virginia

North Carolina

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																	37.7		.sgs.com	
CLIENT:	TEC INC.				SGS Refe	erence #:											page		of	
CONTACT:	Jeff Hart F	HONE NO:	303.273.0231									,		<u>,</u>			,,		·	
PROJECT:	9121-003 s	ITE/PWSID#:	Red Hill BFSF	<u>.</u>	<u>"</u>	Preserv. Used	/k <sup>2</sup>		/ķ <sup>i</sup>	_	HH		<u>/</u>	$\angle$	_	_	_	_/		
REPORTS TO:	Jeff Hart e	mail <u>ishart@tec</u> cc <u>wmcwhitm</u>	cinc.com an@tecinc.co	om	* c o z	SAMPLE TYPE C =	(a	8)		IMS)			!	,						
INVOICE TO:	<del>-</del>	QUOTE #: P.O. NUMBER:			T A 1	COMP G =	3RO (8015B)	TPH-DRO (8015B)	VOC's (8260B)	PAH's (8270C-SIMS)	Diss Pb (6020)				٨		<u> </u>			
LAB, N	SAMPLE IDENTIFICATION	DATE	TIME	MATRIX	R S	GRAB	TPH-GRO	тРН-£	NOC.	PAH's	Diss F			1	25	<u></u>	6		REMARKS	
XXXX	RHMW2254-WG10	10/22/2008	0930	Water	10			X		X	X	$\bigcirc$	(	<b>D</b>	(Z)	K		_	3x Volume sent in 2 co	olers
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				<del>  ,</del>	<u> </u>		<del> </del>	<u> </u>						<u> </u>						
Collected/Relinquish	ned By: (1)	Date	Time	Received By	<del>\</del>	1	17	l	Shippin	a Corri	or:		•				Samples	: Recei	red Cold? YES NO	-
لألا	10. W.	16-22-48	1720		LA	[] 	U_		Shippin								Tempera			
Relinquished By	(2)	Date	Time	Received By:	•	-			Special	Delive	rable F	Requiren	ients:				Chain c	of Custo	dy Seal: (Circle)	
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Relinquished By: (	(3)	Date	Time	Received By:			ich	2	Reques	ted Tu	meroun	đ Time	and-or	Specia	al Instru	ctions:				
						10/211	$I_{i\alpha}$		See											
Relinquished By:	(4)	Date	Time	Received For Lab	oratory By	<i>Y-</i> _	<del>-</del>	•	,									-	_	
200 W. Po	tter Drive Anchorage, AK 99518 Tel: (9	07) 562-2343 Fax: (907)	561-5301	- TU		<u> </u>	James	Drive V	Vest St	Rose,	LA 70	0 <b>087</b> Te	el: (504	) 469- <del>(</del>	6401 Fa	x: (50	4) 463-3	304		

3180 Peger Road Fairbanks, AK 99701 Tel: (907) 474-8656 Fax: (907) 474-9685

255 Sand Island Access Rd., Unit 1B Honolulu, HI 96819 Tel: (808) 224-6217 Fax: (808) 845-2287

1258 Greenbrier Street Charleston, WV 25311 Tel: (304) 346-0725 Fax: (304) 346-0761

5500 Business Drive Wilmington, NC 28405 Tel: (910) 350-1903 Fax: (910) 350-1557





#### Locations Nationwide

Alaska

Hawaii

West Virginia

Maryland

i Louisiana

New Jersey

North Carolina

www.us.sgs.com SGS Reference #: CLIENT: TEC INC. of \_\_\_\_ 303.273.0231 CONTACT: Jeff Hart PHONE NO: 9121-003 Red Hill BFSF Used SITE/PWSID#: PROJECT: SAMPLE jshart@tecinc.com email REPORTS TO: Jeff Hart C wmcwhitman@tecinc.com 0 PAH's (8270C-SIMS) C = TPH-DRO (8015B) TPH-GRO (8015B) COMP QUOTE #: TEC INC INVOICE TO: P.O. NUMBER: 요 G = GRAB SAMPLE IDENTIFICATION TIME MATRIX DATE LAB NO. REMARKS X X 5 X 10/22/2008 1245 Water RHMW02-WG13 X X X Water 5 RHMWA01-WG13 10/22/2008 1205 Received By: Collected/Relinquished By: (1) Samples Received Cold? YES NO Shipping Carrier. Temperature °C:TB=25 C=2.6 Shipping Ticket No: Received By: Relinquished By: Chain of Custody Seal: (Circle) Special Deliverable Requirements: INTACT BROKEN ABSENT See Contract Received By: Relinquished By: (3) Requested Turnaround Time and-or Special Instructions: See Contract Received For Laboratory By Relinguished By: (4) 1050 151 James Drive West St Rose, LA 70087 Tel: (504) 469-6401 Fax: (504) 463-3304 200 W. Potter Drive Anchorage, AK 99518 Tel: (907) 562-2343 Fax: (907) 561-5301 1258 Greenbrier Street Charleston, WV 25311 Tel: (304) 346-0725 Fax: (304) 346-0761 3180 Peger Road Fairbanks, AK 99701 Tel: (907) 474-8656 Fax: (907) 474-9685 5500 Business Drive Wilmington, NC 28405 Tel: (910) 350-1903 Fax: (910) 350-1557 255 Sand Island Access Rd., Unit 1B Honolulu, HI 96819 Tel: (808) 224-6217 Fax: (808) 245-229



1085813

Locations Nationwide

Hawaii

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Louisiana

rsey West Virginia

Carolina

www.us.sgs.com

			<del></del>	····	SGS Refe	rence #:									Т			
CLIENT: CONTACT:	TEC INC.	PHONE NO:	303.273.0231		1										1	pa	ge _	of
PROJECT:	9121-003	<del></del>	Red Hill BFSF			Preserv.	/K <sup>O</sup>		/k <sup>c</sup> i		KHY		$\overline{Z}$	//				
REPORTS TO:	Jeff Hart	email <u>jshart@tec</u> ∞ <u>wmcwhitm</u>	inc.com an@tecinc.co	<u>om</u>	# 0 8	SAMPLE TYPE C =	6	3}		MS)								
INVOICE TO:	TEC INC	QUOTE #: P.O. NUMBER:			† T : A ! N	COMP G =	TPH-GRO (8015B)	TPH-DRO (8015B)	(8260B)	(8270C-SIMS)	Diss Pb (6020)	·						
LAB NO.	SAMPLE IDENTIFICATION	N DATE	TIME	MATRIX	F R S	GRAB	TPH-G	трн-	VOC's	PAH's	Diss P							REMARKS
76-5	RHMW01-WG13	10/22/2008	1500	Water	<b>1</b> 54			Х		X	X							Filter Diss. Pb in Lab
DG-K	RHMW03-WG13	10/22/2008	1115	Water	5_		<u> </u>	X		X	X					<u> </u>		
Collected/Relinquis	D. Wille	Date 10/22/08 Date 10/23/03	Time  /72  Time	Received By:	1	) }	1	)	Shippir	ng Carr	et No:	Requirem	ents:			Tem	/ perature	ceived Cold? YES NO  *c:78=2.0. C=3.5  ustody Seal: (Circle)
ļ		10/23/65	14:30						See (	Contra	ct					INTA	CT E	BROKEN ABSENT
Relinquished By:	(3)	Date	Time	Received By:					Reque	steđ Tu	ırnaroun	d Time	and-or	Special	Instruction	ıs:		
Relinquished By:	(4) Otter Drive Anchorage, AK 99518 Te	Date 1/2 / 7 8 1: (907) 562-2343 Fex: (907)	Time 1050 561-5301	Received For Lab	oratory By:				West Si	t Rose		<b>0087</b> Te		469-640				
☐ 3180 Pege	r Roed Feirbanks, AK 99701 Tel: (9 Island Access Rd., Unit 18 Honolulu,	907) 474-8656 Fax: (907) 474	1-9685	2287										350-190 (350-190				



Form # F004r17 revised 04/11/08

# SAMPLE RECEIPT FORM

SGS WO#:

Yes	No	NA	_	
_0	_		Are samples RUSH, priority or w/in 72 hrs of hold time?	TAT (circle one): Standard or- Rush
			If yes, have you done e-mail ALERT notification?	Received Date: 10-24-08
	<u></u>		Are samples within 24 hrs. of hold time or due date?	Received Time: /050
			If yes, have you also <i>spoken with</i> supervisor?	Is date/time conversion necessary?
			Archiving bottles (if req'd): Are they properly marked?	# of hours to AK Local Time: +2
<u>*                                    </u>	·		Are there any problems? PM Notified?	Thermometer ID: 69D, 70D
F			Were samples preserved correctly and pH verified?	Cooler ID Temp Blank Cooler Temp
			<del></del>	2 2.0°C 2.0°C 5.6°C
			<u> </u>	$\frac{2}{3}$ $\frac{3.1}{2.5}$ $\stackrel{\circ}{\circ}$ $\frac{2.6}{2.6}$ $\stackrel{\circ}{\circ}$
		•	If this is for PWS, provide PWSID.	<u>y</u> <u>2.0 °C 3.5 °C</u>
			Will courier charges apply?	°C °C
			Method of payment?	Note: Temperature readings include thermometer correction factors
<u> </u>			Data package required? (Level: 1 / 2 / 3 / 4)	Delivery method (circle all that apply): Client /
			Notes:	Alert Courier / UPS / FedEx/ USPS / DHL /
			Is this a DoD project? (USACE, Navy, AFCEE)	AA Goldstreak / NAC / ERA / PenAir / Carlile/
22220	Q <u>00</u> 000000	(10.000.000.000.0000.0000.0000.0000.000		Lynden / SGS / Other:
200 X X X X X X X X X X X X X X X X X X			must be filled out for DoD projects (USACE, Navy, AFCEE)	Airbill # 8665 6327 7905 Additional Sample Remarks: (√if applicable)
Yes	ì	o	Is received temperature 4 ± 2°C?	Extra Sample Volume?
	emmer <u>edi</u> Kananaka Kananaka		Exceptions: Samples/Analyses Affected:	Limited Sample Volume?
22222 22222 22222				MeOH field preserved for volatiles?
XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX	********* *********			Field-filtered for dissolved /- 6
222222 222222 222222				Lab-filtered for dissolved
XXXXXXX XXXXXXX	XXXXXXXXX XXXXXXXXXXXXXXXXXXXXXXXXXXXX		If temperature(s) < 0 °C, were containers ice-free? (N/A)  Notify PM immediately of any ice in samples.	Ref Lab required?
		Y O O O O O O O O O O O O O O O O O O O	Was there an airbill? (Note # above in the right hand column)	Foreign Soil?
			Was cooler sealed with custody seals?	
NEXXXX XXXXXX XXXXXX	XUMMANK XXXXXXXX XXXXXXX	COMMANA COMMANA COMMANA	#/where: /on front/on back of a V	This section must be filled if problems are found.  Yes No
~~~		6 N K K K K K K K K 6 K K K K K K K K <del>6 K K K</del> K K K K K K 7 K K K K K K K K K K K K K	Were seal(s) intact upon arrival?	Was client notified of problems?
ين ا		enerenen <del>ener</del> enen enere	Was there a COC with cooler? Was COC sealed in plastic bag & taped inside lid of cooler?	
	<u> </u>		Was the COC filled out properly?	Individual contacted:
******************	********* ********	-	:Did the COC indicate USACE / Navy / AFCEE project?	Via: Phone / Fax / Email (circle one)  Date/Time:
T.			Did the COC and samples correspond?	Reason for contact:
2	<b>/</b>	GRANNEN Granning Granning	Were all sample packed to prevent breakage?	Accepted for contact.
XXXXXXX			Packing material: (300)  Were all samples unbroken and clearly labeled?	
		K	Were all samples end oken and clearly labeled:  Were all samples sealed in separate plastic bags?	
	<u>~~~~~~~</u> ×××××××		Were all VOCs free of headspace and/or MeOH preserved?	
×××××× ××××××	*-		Were correct container / sample sizes submitted?	Change Order Required?
*~			Is sample condition good?	SGS Contact:
			Was copy of CoC, SRF, and custody seals given to PM to fax?	
RERERE		EWEREKK!	3.0	
Note	s: <u>-</u> *	<u>: كأن</u>	ent wants sample 1 filtered at lab bu	ut sample is already
	•			
		177	log preserved to a pH < 2.	
V 6	111	-	VAL RY-CSIC	
外(2	>MH-1	15 19	VOAS (SX-C>1cm	1 3 4 61 01
du	101 6	<i>کم حم</i> د	ed with TB analysis. Hold off on dissor	Wed 76 for sumple #7. 10/246-
400	-0-	no CC I	10000	
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	pleted			- Court Soul
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pg) of 2

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# SAMPLE RECEIPT FORM (page 2)

SGS WO#:

Container							er V	olur.	ne		Container Type							Preservative												
#	Container ID	Matrix	Test	OC	TB	11	500 mL	_	125 mL	60 mL	40 mL	8oz (250 mL)	4oz (125 mL)	Other	AG	ĐO	HDPE	Nalgene	Cubie	Coli	Septa	Other	None	HCI	HNO3	H <sub>2</sub> SO <sub>4</sub>	МеОН	$Na_2S_2O_3$	NaOH	Other
)	A-C	l	TPH-GA-0								3				X						X			X						
	D-F		VOC 8260								3				×						Ż			X						
	G		DissPb TPH-D2-08085					,									$\times$				,			<u> </u>	X					FF
	H-I		TPH-DA-08015			Z									X							-	XX			<u> </u>				
	JK		194 8270 SIMS			2									X								X							
<u> </u>	L-M		extravol	n			2										X				~			2.6	X	ļ				-
2	A-B			ns							2				<b>/</b> _	-	_				X			X.	<u> </u>	<u> </u>			-	
	C-E		Voc DissPb	-							3	5 <i>A</i> (	3		$\succ$		-				X			$\not\succeq$						
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	G-H I-J		TPH-DRO	17		ત્રીત્ર									$\Rightarrow$		-								<del>                                     </del>					
3	4-C	1	PAH-822051MS	ns		<u>~</u>					3				文						$\overline{}$		7	X	<del> </del>				.	· ·
<u></u>	D-F		TPH-G	1						-	3				Ż						45			V	<u> </u>					
	5		Diss Pb									ΓA	1)	,	~									/-	<u> </u>					
	H-I		TPH-DRO			X				ľ		,,,			X									X						
	J-K		PAH8270csus	V		Q								<del></del>	$\checkmark$									/						
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Bottle Totals	123	16	

Completed by: \_

Date: 10-24-0

Form # F004r16 revised 03/10/03

Pg 2012

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# SAMPLE RECEIPT FORM (page 2)

SGS WO#:

	1.4.2.2.2	11.	ere e de la companya	şr .			Container Volume							Container Type								Preservative								
#	Container ID	Matrix	Test	) )	TB	IL	500 mL	250 mL	125 mL		40 mL		4oz (125 mL)	Other	AG	90	HDPE	Nalgene	Cubie	Coli	Septa	Other	None	IDH	HINO3	$\mathrm{H}_2\mathrm{SO}_4$	МеОН	$Na_2S_2O_3$	NaOH	Other
4-6	A-c	1	TPH-GRO								0				Χ						Y			X						
	D-F		VOC 8260				G	7-6	6		9	·			X						$\times$			X					$\square$	
	G H-I		Diss Plo TPH-DRO8015				1	2			· · · · · · · · · · · · · · · · · · ·						X				<u>'</u>				X					FE
<u> </u>						6					<u></u>				オオ									X					$\square$	
	J-K		PAH 8270c S(M)			6							-		X								X						$\vdash$	
7	A-C	•	TPH GRD 8015/3								マ				\ <u>\</u>						X		$\vdash$	V						
	D-F	1	VOC8260								3				X				-		$\langle \cdot \rangle$			1	_				*	
	4		DissPb					٢			<u> </u>				$\rightarrow$		X				^-									
	H-I		Diss Pb TPH-DROGOLS			2	<del></del> -—-	.1		_					X									X						
	G H-エ ナ		PAH 82705 SIMS												X								X							
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Bottle Totals 15 1 3 27

Completed by: \_

-Date: 10-24-68

Form # F004r16 revised 03/10/03

0	Fedex US Airbill 8665 6327 790.	5 0200 for FedEx Retrieval Copy
0	11 From Date 102308 Sender's FedEx Account Number	4a Express Package Service  1 FedEx Priority Overnight Next business morning. Fridey Shipmens will be delivered on Monaday shipmens will be delivered on Monaday shipmens will be delivered on Monaday shipmens will be delivered in Monaday shipmens will be delive
0	Sender's Name Phone CD F47 (10) 6 )	3   Second business day. Thursday   20   The dex Express Saver through the delevated on Monday unless SATURDAY Cellsons a selection 12   Second business satisfaction of the selection of the s
1800.463.3339	Company ESM PACIFIC  Address 1510 Kayan 27.  Dept. Macrobine Roo.	7 FedEx 1Day Freight* Next business day: The State of the
	city Huriolain State HI ZIP 96818	FedEx 2 1 Albert public small pake. FedEx 4 FedEx 1 Other Tube. FedEx large Pak and FedEx South
O 1.800.GoFedEx	2 Your Internal Billing Reference  3 To Recipient's Name Phone 7-7 5-12 2344	Envelope*  In the Decired value limit \$500.  Special Handlin  Special Handlin  Include FedEx By in Section 3.  SATURDAY Delivery Covernment of Fedex Location 3.  Saturday Delivery Covernment of Fedex Delivery C
C dedex.com	Company 363 ENXILAINENTER DERVICES	Does this shipment contain dangerous goods?  One box must be checked.  Yes As per attached As per attached to the shipper of the box to the shipper of the shipper of the box to the shipper of the box
	Recipient's 200 W. Ps 77 CC DR.  We cannot deliver to P.O. boxes or P.O. ZIP codes.  Address  Address	7 Payment Bill to: Enter FedEx Acet. No. or Credit Card No. below. 1 Sender Acet No. no Credit Card No. below. 2 Recipient 3 Third Party 4 Credit Card 5 Cash/Check
0	To request a package be held at a specific FedEx location, print FedEx address here.  City ACAMMA State State AK ZIP 97518	Total Packages Total Weight
0		Our liability is limited to \$100 unless you declare a higher value. See the current Fedex Service Guide for dotals. Credit Card Auth:  8 Residential Delivery Signature Options If you require a signature, check Director Indirect.
		No Signature  Required  Parking may be left, 10 10 diddress may sign for distance, so someone as a neighboring address any sign for distance, so someone at a neighboring address nay sign for distance, for degrees and some as a neighboring address nay sign for distance, for degrees and some as a neighboring address nay sign for distance, for degrees, for degrees and some



# SGS Environmental Services Alaska Division Level II Laboratory Data Report

Project:	9121-003 Red Hill BFSF

Client: The Environmental Company, Inc. (TEC)

SGS Work Order: 1086126

Released by:

#### Contents:

Cover Page
Case Narrative
Final Report Pages
Quality Control Summary Forms
Chain of Custody/Sample Receipt Forms

#### Note:

Unless otherwise noted, all quality assurance/quality control criteria is in compliance with the standards set forth by the proper regulatory authority, the SGS Quality Assurance Program Plan, and the National Environmental Accreditation Conference.



CASE NARRATIVE Print Date: 12/11/2008

Client Name: The Environmental Company, Inc. (TEC)

Project Name: 9121-003 Red Hill BFSF

Workorder No.: 1086126

# Sample Comments

Refer to the sample receipt form for information on sample condition.

<u>Lab Sample ID</u> <u>Sample Type</u> <u>Client Sample ID</u>

There were no analytical anomalies associated with the data reported herein.



# Laboratory Analytical Report

Client: The Environmental Company, Inc.

1001 Bishop Street Ste 1400

**ASB Tower** 

Honolulu, HI 96813

Attn: Jeff Hart

T: (808)528-1445 F:(808)528-0768

jshart@tecinc.com

Project: 9121-003 Red Hill BFSF

Workorder No.: 1086126

#### Certification:

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, other than the conditions noted on the sample data sheet(s) and/or the case narrative. This certification applies only to the tested parameters and the specific sample(s) received at the laboratory.

If you have any questions regarding this report, or if we can be of further assistance, please contact your SGS Project Manager.

Tamara Rentz tamara.rentz@sgs.com Project Manager



Print Date: 12/11/2008

Enclosed are the analytical results associated with this workorder.

As required by the state of Alaska and the USEPA, a formal Quality Assurance/Quality Control Program is maintained by SGS. A copy of our Quality Assurance Plan (QAP), which outlines this program is available at your request.

The Laboratory certification numbers are AK971-05 (DW), UTS-005 (CS) and AK00971 (Micro) for ADEC and AK100001 for NELAP (RCRA methods: 1020A, 1311, 6010B, 7470A, 7471A, 9040B, 9045C, 9056, 9060, 8015B, 8021B, 8081A/8082, 8260B, 8270C).

Except as specifically noted, all statements and data in this report are in conformance to the provisions set forth by the SGS QAP, the National Environmental Laboratory Accreditation Program and, when applicable, other regulatory authorities.

If you have any questions regarding this report or if we can be of any assistance, please contact your SGS Project Manager at 907-562-2343.

The following descriptors may be found on your report which will serve to further qualify the data.

MDL Method Detection Limit

PQL Practical Quantitation Limit (reporting limit).

CL Control Limit

U Indicates the analyte was analyzed for but not detected. F Indicates value that is greater than or equal to the MDL.

J The quantitation is an estimation.

ND Indicates the analyte is not detected

B Indicates the analyte is found in a blank associated with the sample.

\* The analyte has exceeded allowable regulatory or control limits.

D The analyte concentration is the result of dilution.

GT Greater Than LT Less Than

Q QC parameter out of acceptance range.

M A matrix effect was present.

E The analyte result is above the calibrated range.

R Rejected

DF Analytical Dilution Factor

JL The analyte was positively identified, but the quantitation is a low estimation.

<Surr> Surrogate QC spiked standard

<Surr/IS> Surrogate / Internal Standard QC spiked standard

QC Quality Control
QA Quality Assurance
MB Method Blank

LCS (D) Laboratory Control Sample (Duplicate)

MS(D) Matrix Spike (Duplicate)
BMS(D) Site Specific Matrix Spike
RPD Relative Percent Difference
ICV Initial Calibration Verification
CCV Continuous Calibration Verification
MSA Method of Standard Addition

Notes: Soil samples are reported on a dry weight basis unless otherwise specified

All DRO/RRO analysese are integrated per SOP.



SAMPLE SUMMARY

Print Date: 12/11/2008 1:06 pm

Client Name: The Environmental Company, Inc. (TEC)

Project Name: 9121-003 Red Hill BFSF

Workorder No.: 1086126

**Analytical Methods** 

**Analytical Method** Method Description

Dissolved Metals by ICP-MS SW6020

Sample ID Cross Reference

Lab Sample ID Client Sample ID 1086126001 RHMW01-WG13



## The Environmental Company, Inc. (TEC)

Client Sample ID: RHMW01-WG13

SGS Ref. #: 1086126001

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 11/03/08 13:35 Receipt Date/Time: 11/05/08 09:45

Print Date: 12/11/2008 1:06 pm

## Dissolved Metals by ICP/MS

<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Analytical Batch	<u>Prep</u> Batch	Qualifiers		
Lead	0.966 J	1.00	0.310	ug/L	5	MMS5721	MXX21078	8		
Batch Information										
Analytical Batch: MMS5721		Prep Batch	: MXX21078			Initial Prep	Wt./Vol.: 50 ı	mL		
Analytical Method: SW6020		Prep Metho	d: SW3010A			Prep Extra	Prep Extract Vol.: 50 mL			
Analysis Date/Time: 11/15/08 17:30		Prep Date/	Γime: 11/13/08 2	0:00		Container I	D:108612600	01-A		
Dilution Factor: 5						Analyst: BN	ИΕ			



SGS Ref.# Client Name 871318

Method Blank

The Environmental Company, Inc. (TEC)

Project Name/# Matrix 9121-003 Red Hill BFSF

Water (Surface, Eff., Ground)

Printed Date/Time

Prep

12/11/2008 13:06

Batch Method

MXX21078 SW3010A

Date 11/13/2008

QC results affect the following production samples:

1086126001

Parameter		Results	Reporting/Control Limit	MDL	Units	Analysis Date
Metals by ICI	P/MS					
Lead		ND	1.00	0.310	ug/L	11/20/08
Batch	MMS5732					
Method	SW6020					
Instrument	Perkin Elmer Sciex ICP-MS F	24				



SGS Ref.# 871319 Lab Control Sample Printed Date/Time

Prep Batch MXX

12/11/2008

13:06

Batch MXX21078 Method SW3010A

**Date** 11/13/2008

Client Name The Environmental Company, Inc. (TEC)
Project Name/# 9121-003 Red Hill BFSF

Matrix Water (Surface, Eff., Ground)

QC results affect the following production samples:

1086126001

Parameter QC Pct LCS/LCSD RPD Spiked Analysis
Parameter Results Recov Limits RPD Limits Amount Date

Metals by ICP/MS

Lead LCS 1000 100 (80-120) 1000 ug/L 11/20/2008

Batch MMS5732 Method SW6020

**Instrument** Perkin Elmer Sciex ICP-MS P4



# CHAIN OF CUSTODY RECORD SGS Environmental Services Inc.

1086126

.ocations Nationwide

Hawaii

Louisiana

West Virginia

ersey Carolina

															عبعميس	<u>w</u>	ww.u	s.sgs.com
CLIENT:	TEC INC.				SGS Ref	erence #:										page		of
CONTACT:	Jeff Hart	PHONE NO:	303.273.0231													P-50		
PROJECT:	9121-003	SITE/PWSID#:	Red Hill BFSF	,		Preserv.	Į,		K.C.		KHIC		//	//	/		_	
REPORTS TO:	Jeff Hart	email jshart@te	ecinc.com nan@tecinc.co	o <u>m</u>		SAMPLE TYPE C =	<b>6</b>	e e		MS)								
INVOICE TO:	TEC INC	OUOTE #: P.O. NUMBER:			T A I N	COMP G·=	TPH-GRO (8015B)	TPH-DRO (8015B)	(8260B)	(8270C-SIMS)	Pb (6020)						·	
LAB NO.	SAMPLE IDENTIFICATION	N DATE	TIME	MATRIX	E FI S	GRAB	TPH-G	TPH-C	Voc's	PAH's	Diss F							REMARKS
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Collected/Relinquish	20. W.L.	Date 11/4/09	Timew 094	Received By:	<del></del>				l	g Carri g Ticke		_ · ·	•	•	•	Samples 690 Tempera	Rece <b>2</b> ature	ived Cold? YES NO FB=1.4C = 1.0
elinquished By: (	2)	Date	Time	Received By:						Delive		equireme	nts:		•	Chain o		ody Seal: (Circle)
Relinquished By: (	3)	Date	Tilite	Received By:					Reques	ited Tu	rnaround	l Time a	ınd-or Sp	ecial Insti	ructions			
									See	Con	itrac	į						
Relinquished By: (	4)	11/5/08	Time 1//0	Received for Labe	yatory By:													
3180 Peger	tter Drive Anchorage, AK 99518 Tel: Road Fairbanks, AK 99701 Tel: (90 sland Access Rd., Unit 18 Honolutu,	(907) 562-2343 Fax: (907) 07) 474-8656 Fax: (907) 47	4-9685	<del></del>		151 1258 5500	Greenb	rier St	reet Ch	ariestor	, WV	25311 T	el: (304)	346-0725	Fax:	(304) 34	6-0761	

## 1086126



## SAMPLE RECEIPT FORM

SGS WO#:

Yes	No	NA		
1 40	1	14/7	Are samples RUSH, priority or w/in 72 hrs of hold time?	TAT (circle one): Standard -or- Rush
	<u> </u>	7-	If yes, have you done <i>e-mail ALERT notification</i> ?	Received Date: 1/57-08
			Are samples within 24 hrs. of hold time or due date?	Received Time: ///O
<del></del> .	1	1	If yes, have you also <i>spoken with</i> supervisor?	Is date/time conversion necessary? WO
	—		Archiving bottles (if req'd): Are they properly marked?	# of hours to AK Local Time:
				Thermometer ID: 696
			Are there any problems? PM Notified?	
			Were samples preserved correctly and pH verified?	Cooler ID Temp Blank Cooler Temp
			<del></del>	· · · · · · · · · · · · · · · · · · ·
		1_	If this is for PWS, provide PWSID.	
			Will courier charges apply?	
			Method of payment?	
			Data package required? (Level: 1 / 2 / 3 / 4)	Note: Temperature readings include thermometer correction factors <b>Delivery method</b> (circle all that apply): Client /
			Notes:	Alert Courier / UPS FedEy / USPS / DHL /
			is this a DoD project? (USACE, Navy, AFCEE)	AA Goldstreak / NAC / ERA / PenAir / Carlile/
—			is this a bob project: (boxot, navy, ra ozz,	Lynden / SGS / Other:
2202233	ri.		must be filled out for DoD projects (USACE, Navy, AFCEE)	Airbill # 7911 8073 0793
Yes		o cuon	made of pateu out for Door projects (ODACE, May), ATCEE]	Additional Sample Remarks: (\(\sqrt{if applicable}\)
XXXXXXXX XXXXXXXX XXXXXXX		<i>_</i>	Is received temperature 4 ± 2°C?	Extra Sample Volume?
	växk <del>sa</del>		Exceptions: Samples/Analyses Affected:	Limited Sample Volume?
		6 2 3 2 2 2 3 3 2 2 4 3 3 2 2 3 3 3 3 3 4 3 3 3 3 3 3 3 3 3 3 3 3 3 3	SEE ABOVE	MeOH field preserved for volatiles?
KERKERN KERKEKK KERKEKK	XXXXXXXX XXXXXXXX XXXXXXX			Field-filtered for dissolved %
2000 2000 2000 2000 2000				Lab-filtered for dissolved
202020 20 <del>4040</del>	N	A	If temperature(s) <0 °C, were containers ice-free? N/A	Ref Lab required?
			Notify PM immediately of any ice in samples.  Was there an airbill? (Note # above in the right hand column)	Foreign Soil?
4			Was cooler sealed with custody seals?	
	XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX		#/ where: 2 FRONT & BJOK TOP UP	This section must be filled if problems are found.
سا		XXXXXXXX XXXXXXXX	Were seal(s) intact upon arrival?	Yes No
س	**************************************		Was there a COC with cooler?	Was client notified of problems?
س)			Was COC sealed in plastic bag & taped inside lid of cooler?	Individual contacted:
- U - U - U	Yakkari Takkari Kan		Was the COC filled out properly?	Via: Phone / Fax / Email (circle one)
-	<del>Ma</del> nnan Xannan Yannan	KMMXMAK KKKKKKK TOKKKKK	Did the COC indicate USACE / Navy / AFCEE project?	Date/Time:
***********			Did the COC and samples correspond?	Reason for contact;
	ANNANA Kanana	SKEKSEK <del>TOTO</del> LEKE	Were all sample packed to prevent breakage?	
	XXXXXX		Packing material: Bospus wydd Were all samples unbroken and clearly labeled?	
×××××××		<u> </u>	Were all samples sealed in separate plastic bags?	
~	· · · · · · · · · · · · · · · · · · ·	/A-	Were all VOCs free of headspace and/or MeOH preserved?	
V			Were correct container / sample sizes submitted?	
v			Is sample condition good?	Change Order Required?
L	XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX		Was copy of CoC, SRF, and custody seals given to PM to fax?	SGS Contact:
KANAMAK KANAMAK KANAMAK	RMKRMK RMKRMK			
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Notes	»:			
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## SAMPLE RECEIPT FORM (page 2)

SGS WO#:

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#		L I	Test				,							Other		T						Other								Other
17	Container ID	Matrix	1051	) ()	TB	.1L	500 mL	250 mL	125 mL	60 mL	40 mL	8oz (250 mL)	4oz (125 mL)	Other	AG	90	HDPE	Nalgene	Cubie	Coli	Septa	Other	None	HCI	HINO3	H <sub>2</sub> SO <sub>4</sub>	МеОН	$Na_2S_2O_3$	NaOH	Other
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Bottle Totals

Completed by:

Date: 1/-5-08

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1086126

From: Origin ID: HIKA (808)528-1445 BILL WHITMAN TEC INC. 1001 BISHOP STREET, ASB TOWER

**SUITE 1400** HONOLULU, HI 96813

SHIP TQ: 9075622343

**BILL THIRD PARTY** 

SAMPLE RECEIVING SGS Environmental Services 200 W POTTER DR

**ANCHORAGE, AK 995181605** 



CAD: 1774997/INET8091 Accoun#; S \*\*\*\*\*\*\*\* Delivery Address Bar Code

> Invoice # PO#

Dept#

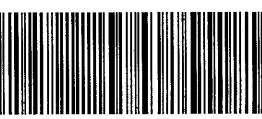
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WED - 05NOV AM PRIORITY OVERNIGHT

WU CYMA

99518 AK-US **ANC** 



After printing this label:

1. Use the 'Print' button on this page to print your label to your laser or inkjet printer.

2. Fold the printed page along the horizontal line.

3. Place label in shipping pouch and affix it to your shipment so that the barcode portion of the label can be read and scanned.

Warning: Use only the printed original label for shipping. Using a photocopy of this label for shipping purposes is fraudulent and could result in additional billing charges, along with the cancellation of your FedEx account number.

Use of this system constitutes your agreement to the service conditions in the current FedEx Service Guide, available on fedex.com.FedEx will not be responsible for any claim in excess of \$100 per package, whether the result of loss, damage, delay, non-delivery, misdelivery, or misinformation, unless you declare a higher value, pay an additional charge, document your actual loss and file a timely claim. Limitations found in the current FedEx Service Guide apply. Your right to recover from FedEx for any loss, including intrinsic value of the package, loss of sales, income interest, profit, attorney's fees, costs, and other forms of damage whether direct, incidental, consequential, or special is limited to the greater of \$100 or the authorized declared value. Recovery cannot exceed actual documented loss. Maximum for items of extraordinary value is \$500, e.g. jewelry, precious metals, negotiable instruments and other items listed in our ServiceGuide. Written claims must be filed within strict time limits, see current FedEx Service Guide.



# SGS Environmental Services Alaska Division Level II Laboratory Data Report

Project:	9121-003 Red Hill BFSF

Client: The Environmental Company, Inc. (TEC)

SGS Work Order: 1086721

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#### Contents:

Cover Page
Case Narrative
Final Report Pages
Quality Control Summary Forms
Chain of Custody/Sample Receipt Forms

#### Note:

Unless otherwise noted, all quality assurance/quality control criteria is in compliance with the standards set forth by the proper regulatory authority, the SGS Quality Assurance Program Plan, and the National Environmental Accreditation Conference.



Print Date: 12/22/2008 **CASE NARRATIVE** 

Client Name: The Environmental Company, Inc. (TEC)

Project Name: 9121-003 Red Hill BFSF

Workorder No.: 1086721

#### Sample Comments

Refer to the sample receipt form for information on sample condition.

Lab Sample ID	Sample Type	Client Sample ID

876591 MS 1086597003A(876590MS)

8260B - MS recovery for several analytes does not meet QC goals (biased high). See LCS for accuracy.

876592 MSD 1086597003A(876590MSD)

8260B - MSD recovery for several analytes does not meet QC goals (biased high). See LCS for accuracy.

876594 CCV for HBN 210083 [VMS/10317]

8260B - ICV recovery for dichlorodifluoromethane and 1,1-dichloroethene does not meet QC goals (biased high). These

analytes were not detected above the PQL in the associated samples.



## Laboratory Analytical Report

Client: The Environmental Company, Inc.

1001 Bishop Street Ste 1400

**ASB Tower** 

Honolulu, HI 96813

Attn: Jeff Hart

T: (808)528-1445 F:(808)528-0768

jshart@tecinc.com

Project: 9121-003 Red Hill BFSF

Workorder No.: 1086721

#### Certification:

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, other than the conditions noted on the sample data sheet(s) and/or the case narrative. This certification applies only to the tested parameters and the specific sample(s) received at the laboratory.

If you have any questions regarding this report, or if we can be of further assistance, please contact your SGS Project Manager.

Tamara Rentz

tamara.rentz@sgs.com

Project Manager



Print Date: 12/22/2008

Enclosed are the analytical results associated with this workorder.

As required by the state of Alaska and the USEPA, a formal Quality Assurance/Quality Control Program is maintained by SGS. A copy of our Quality Assurance Plan (QAP), which outlines this program is available at your request.

The Laboratory certification numbers are AK971-05 (DW), UTS-005 (CS) and AK00971 (Micro) for ADEC and AK100001 for NELAP (RCRA methods: 1020A, 1311, 6010B, 7470A, 7471A, 9040B, 9045C, 9056, 9060, 8015B, 8021B, 8081A/8082, 8260B, 8270C).

Except as specifically noted, all statements and data in this report are in conformance to the provisions set forth by the SGS QAP, the National Environmental Laboratory Accreditation Program and, when applicable, other regulatory authorities.

If you have any questions regarding this report or if we can be of any assistance, please contact your SGS Project Manager at 907-562-2343.

The following descriptors may be found on your report which will serve to further qualify the data.

MDL Method Detection Limit

**PQL** Practical Quantitation Limit (reporting limit).

CL Control Limit

П Indicates the analyte was analyzed for but not detected. F Indicates value that is greater than or equal to the MDL.

J The quantitation is an estimation. ND Indicates the analyte is not detected

В Indicates the analyte is found in a blank associated with the sample. The analyte has exceeded allowable regulatory or control limits.

D The analyte concentration is the result of dilution.

GT Greater Than LT Less Than

Q QC parameter out of acceptance range.

A matrix effect was present. Μ

Ε The analyte result is above the calibrated range.

R Rejected

DF Analytical Dilution Factor

JL The analyte was positively identified, but the quantitation is a low estimation.

<Surr> Surrogate QC spiked standard

<Surr/IS> Surrogate / Internal Standard QC spiked standard

Quality Control QC Quality Assurance QA MB Method Blank

LCS (D) Laboratory Control Sample (Duplicate)

Matrix Spike (Duplicate) MS(D) Site Specific Matrix Spike BMS(D) **RPD** Relative Percent Difference Initial Calibration Verification **ICV** CCV Continuous Calibration Verification MSA Method of Standard Addition

Notes: Soil samples are reported on a dry weight basis unless otherwise specified

All DRO/RRO analysese are integrated per SOP.



SAMPLE SUMMARY

Print Date: 12/22/2008 11:45 am

Client Name: The Environmental Company, Inc. (TEC)

Project Name: 9121-003 Red Hill BFSF

Workorder No.: 1086721

**Analytical Methods** 

Method DescriptionAnalytical MethodAFCEE 3.1 8260 (W)SW8260B

Sample ID Cross Reference

 Lab Sample ID
 Client Sample ID

 1086721001
 RHMW2254-WG13B

 1086721002
 RHMWA01-WG13B

 1086721003
 TB01-WG13B



Client Sample ID: RHMW2254-WG13B

SGS Ref. #: 1086721001

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 12/16/08 10:05 Receipt Date/Time: 12/17/08 11:55

Print Date: 12/22/2008 11:45 am

#### **Volatile Gas Chromatography/Mass Spectroscopy**

	.,,					Analytical	Prep
<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Benzene	ND	0.400	0.120	ug/L	1	VMS10317	VXX19107
Toluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Ethylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
n-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,2-Dichloroethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
4-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Chlorobenzene	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	1	VMS10317	VXX19107
cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
4-Isopropyltoluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
n-Propylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Styrene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Dibromomethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,4-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Acetone	ND	10.0	3.10	ug/L	1	VMS10317	VXX19107
1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,2-Dibromo-3-chloropropane	ND	2.00	0.620	ug/L	1	VMS10317	VXX19107
Methyl-t-butyl ether	ND	5.00	1.50	ug/L	1	VMS10317	VXX19107
Tetrachloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Dibromochloromethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,3-Dichloropropane	ND	0.400	0.120	ug/L	1	VMS10317	VXX19107
1,2-Dibromoethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Carbon tetrachloride	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1,1,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
Chloroform	ND	1.00	0.300	ug/L	1	VMS10317	VXX19107
Bromobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,3-Trichloropropane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Chloromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Bromomethane	ND	3.00	0.940	ug/L	1	VMS10317	VXX19107
Bromochloromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Vinyl chloride	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Dichlorodifluoromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107

SGS Environmental Services Inc. Alaska Division 200 West Potter Drive Anchorage Alaska 99518

† (907) 562-2343 † (907) 561-5301 www.us.sgs.com



Client Sample ID: RHMW2254-WG13B

SGS Ref. #: 1086721001

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground)

Collection Date/Time: 12/16/08 10:05 Receipt Date/Time: 12/17/08 11:55 Print Date: 12/22/2008 11:45 am

#### **Volatile Gas Chromatography/Mass Spectroscopy**

<b>C</b> . ,						<u>Analytical</u>	<u>Prep</u>
<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	<u>Batch</u>	Batch Qualifiers
Chloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
sec-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Bromodichloromethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,1-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
2-Butanone (MEK)	ND	10.0	3.10	ug/L	1	VMS10317	VXX19107
Methylene chloride	ND	5.00	1.00	ug/L	1	VMS10317	VXX19107
Trichlorofluoromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
P & M -Xylene	ND	2.00	0.620	ug/L	1	VMS10317	VXX19107
Naphthalene	ND	2.00	0.620	ug/L	1	VMS10317	VXX19107
o-Xylene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Bromoform	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1-Chlorohexane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
tert-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1-Dichloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
2-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Trichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
2,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Hexachlorobutadiene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Isopropylbenzene (Cumene)	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2-Dichloroethane-D4 <surr></surr>	99.2	73-120		%	1	VMS10317	VXX19107
Toluene-d8 <surr></surr>	99.6	80-120		%	1	VMS10317	VXX19107
4-Bromofluorobenzene <surr></surr>	100	76-120		%	1	VMS10317	VXX19107

#### **Batch Information**

Analytical Batch: VMS10317 Prep Batch: VXX19107 Initial Prep Wt./Vol.: 5 mL
Analytical Method: SW8260B Prep Method: SW5030B Prep Extract Vol.: 5 mL
Analysis Date/Time: 12/19/08 02:19 Prep Date/Time: 12/18/08 14:04 Container ID:1086721001-A

Dilution Factor: 1 Analyst: DSH



Client Sample ID: RHMWA01-WG13B

SGS Ref. #: 1086721002

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 12/16/08 12:05 Receipt Date/Time: 12/17/08 11:55

Print Date: 12/22/2008 11:45 am

#### **Volatile Gas Chromatography/Mass Spectroscopy**

	,					Analytical	<u>Prep</u>
<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Benzene	ND	0.400	0.120	ug/L	1	VMS10317	VXX19107
Toluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Ethylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
n-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,2-Dichloroethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
4-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Chlorobenzene	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	1	VMS10317	VXX19107
cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
4-Isopropyltoluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
n-Propylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Styrene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Dibromomethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,4-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Acetone	ND	10.0	3.10	ug/L	1	VMS10317	VXX19107
1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,2-Dibromo-3-chloropropane	ND	2.00	0.620	ug/L	1	VMS10317	VXX19107
Methyl-t-butyl ether	ND	5.00	1.50	ug/L	1	VMS10317	VXX19107
Tetrachloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Dibromochloromethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,3-Dichloropropane	ND	0.400	0.120	ug/L	1	VMS10317	VXX19107
1,2-Dibromoethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Carbon tetrachloride	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1,1,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
Chloroform	ND	1.00	0.300	ug/L	1	VMS10317	VXX19107
Bromobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Chloromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,3-Trichloropropane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Bromomethane	ND	3.00	0.940	ug/L	1	VMS10317	VXX19107
Bromochloromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Vinyl chloride	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Dichlorodifluoromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107

SGS Environmental Services Inc. Alaska Division 200 West Potter Drive Anchorage Alaska 99518

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Client Sample ID: RHMWA01-WG13B

SGS Ref. #: 1086721002

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground)

Collection Date/Time: 12/16/08 12:05 Receipt Date/Time: 12/17/08 11:55 Print Date: 12/22/2008 11:45 am

Prep

Analytical

#### **Volatile Gas Chromatography/Mass Spectroscopy**

Davamatav	Result	PQL/CL	MDI	Unita	DE	Analytical	Prep Ouglifie	
Parameter Chlorooth and			MDL 0.240	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifie	ers
Chloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
sec-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
Bromodichloromethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107	
1,1-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
2-Butanone (MEK)	ND	10.0	3.10	ug/L	1	VMS10317	VXX19107	
Methylene chloride	ND	5.00	1.00	ug/L	1	VMS10317	VXX19107	
Trichlorofluoromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
P & M -Xylene	ND	2.00	0.620	ug/L	1	VMS10317	VXX19107	
Naphthalene	ND	2.00	0.620	ug/L	1	VMS10317	VXX19107	
o-Xylene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
Bromoform	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1-Chlorohexane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
tert-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,1-Dichloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
2-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
Trichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
2,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
Hexachlorobutadiene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
Isopropylbenzene (Cumene)	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,1-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107	
1,2-Dichloroethane-D4 <surr></surr>	95.9	73-120		%	1	VMS10317	VXX19107	
Toluene-d8 <surr></surr>	102	80-120		%	1	VMS10317	VXX19107	
4-Bromofluorobenzene <surr></surr>	99.6	76-120		%	1	VMS10317	VXX19107	

#### **Batch Information**

Analytical Batch: VMS10317 Prep Batch: VXX19107 Initial Prep Wt./Vol.: 5 mL
Analytical Method: SW8260B Prep Method: SW5030B Prep Extract Vol.: 5 mL
Analysis Date/Time: 12/19/08 02:52 Prep Date/Time: 12/18/08 14:04 Container ID:1086721002-A

Dilution Factor: 1 Analyst: DSH



Client Sample ID: TB01-WG13B

SGS Ref. #: 1086721003

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 12/16/08 08:05 Receipt Date/Time: 12/17/08 11:55

Print Date: 12/22/2008 11:45 am

#### **Volatile Gas Chromatography/Mass Spectroscopy**

	,					Analytical	<u>Prep</u>
<u>Parameter</u>	Result	PQL/CL	MDL	<u>Units</u>	<u>DF</u>	Batch	Batch Qualifiers
Benzene	ND	0.400	0.120	ug/L	1	VMS10317	VXX19107
Toluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Ethylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
n-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,2-Dichloroethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,3,5-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
4-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Chlorobenzene	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	1	VMS10317	VXX19107
cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
4-Isopropyltoluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
n-Propylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Styrene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Dibromomethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,4-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Acetone	ND	10.0	3.10	ug/L	1	VMS10317	VXX19107
1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,2-Dibromo-3-chloropropane	ND	2.00	0.620	ug/L	1	VMS10317	VXX19107
Methyl-t-butyl ether	ND	5.00	1.50	ug/L	1	VMS10317	VXX19107
Tetrachloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Dibromochloromethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,3-Dichloropropane	ND	0.400	0.120	ug/L	1	VMS10317	VXX19107
1,2-Dibromoethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Carbon tetrachloride	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1,1,2-Tetrachloroethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
Chloroform	ND	1.00	0.300	ug/L	1	VMS10317	VXX19107
Bromobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Chloromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,3-Trichloropropane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Bromomethane	ND	3.00	0.940	ug/L	1	VMS10317	VXX19107
Bromochloromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Vinyl chloride	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Dichlorodifluoromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107

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Client Sample ID: TB01-WG13B

SGS Ref. #: 1086721003

Project ID: 9121-003 Red Hill BFSF Matrix: Water (Surface, Eff., Ground) Collection Date/Time: 12/16/08 08:05 Receipt Date/Time: 12/17/08 11:55

Print Date: 12/22/2008 11:45 am

#### **Volatile Gas Chromatography/Mass Spectroscopy**

						<u>Analytical</u>	<u>Prep</u>
<u>Parameter</u>	Result	PQL/CL	<u>MDL</u>	<u>Units</u>	<u>DF</u>	<u>Batch</u>	Batch Qualifiers
Chloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
sec-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Bromodichloromethane	ND	0.500	0.150	ug/L	1	VMS10317	VXX19107
1,1-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
2-Butanone (MEK)	ND	10.0	3.10	ug/L	1	VMS10317	VXX19107
Methylene chloride	ND	5.00	1.00	ug/L	1	VMS10317	VXX19107
Trichlorofluoromethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
P & M -Xylene	ND	2.00	0.620	ug/L	1	VMS10317	VXX19107
Naphthalene	ND	2.00	0.620	ug/L	1	VMS10317	VXX19107
o-Xylene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Bromoform	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1-Chlorohexane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,4-Trimethylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
tert-Butylbenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1,1-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1-Dichloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
2-Chlorotoluene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Trichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
trans-1,2-Dichloroethene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
2,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Hexachlorobutadiene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
Isopropylbenzene (Cumene)	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2-Dichloropropane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1-Dichloropropene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,1,2-Trichloroethane	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	1	VMS10317	VXX19107
1,2-Dichloroethane-D4 <surr></surr>	114	73-120		%	1	VMS10317	VXX19107
Toluene-d8 <surr></surr>	99.2	80-120		%	1	VMS10317	VXX19107
4-Bromofluorobenzene <surr></surr>	103	76-120		%	1	VMS10317	VXX19107

#### **Batch Information**

Prep Batch: VXX19107 Initial Prep Wt./Vol.: 5 mL Analytical Batch: VMS10317 Analytical Method: SW8260B Prep Method: SW5030B Prep Extract Vol.: 5 mL Analysis Date/Time: 12/18/08 20:43 Prep Date/Time: 12/18/08 14:04 Container ID:1086721003-A

Dilution Factor: 1 Analyst: DSH



SGS Ref.# Client Name 876587 Method Blank

The Environmental Company, Inc. (TEC)

Project Name/#

9121-003 Red Hill BFSF

Matrix Water (Surface, Eff., Ground)

**Printed Date/Time** 

Prep

12/22/2008 11:45

Batch Method Date

VXX19107 SW5030B 12/18/2008

QC results affect the following production samples: 1086721001, 1086721002, 1086721003

Parameter	Reporting/Control Results Limit MDL Units		Units	Analysis Date	
Volatile Gas Chromatography/Ma	ss Spectro	oscopy			
Benzene	ND	0.400	0.120	ug/L	12/18/08
Toluene	ND	1.00	0.120	ug/L ug/L	12/18/08
	ND ND	1.00	0.310		12/18/08
Ethylbenzene	ND ND	1.00		ug/L	12/18/08
n-Butylbenzene		0.500	0.310	ug/L	12/18/08
1,4-Dichlorobenzene	ND	0.500	0.150	ug/L	12/18/08
1,2-Dichloroethane	ND	1.00	0.150	ug/L	12/18/08
1,3,5-Trimethylbenzene	ND		0.310	ug/L	12/18/08
4-Chlorotoluene	ND	1.00	0.310	ug/L	12/18/08
Chlorobenzene	ND	0.500	0.150	ug/L	
4-Methyl-2-pentanone (MIBK)	ND	10.0	3.10	ug/L	12/18/08
cis-1,2-Dichloroethene	ND	1.00	0.310	ug/L	12/18/08
4-Isopropyltoluene	ND	1.00	0.310	ug/L	12/18/08
cis-1,3-Dichloropropene	ND	0.500	0.150	ug/L	12/18/08
n-Propylbenzene	ND	1.00	0.310	ug/L	12/18/08
Styrene	ND	1.00	0.310	ug/L	12/18/08
Dibromomethane	ND	1.00	0.310	ug/L	12/18/08
trans-1,3-Dichloropropene	ND	1.00	0.310	ug/L	12/18/08
1,2,4-Trichlorobenzene	ND	1.00	0.310	ug/L	12/18/08
Acetone	ND	10.0	3.10	ug/L	12/18/08
1,1,2,2-Tetrachloroethane	ND	0.500	0.150	ug/L	12/18/08
1,2-Dibromo-3-chloropropane	ND	2.00	0.620	ug/L	12/18/08
Methyl-t-butyl ether	ND	5.00	1.50	ug/L	12/18/08
Tetrachloroethene	ND	1.00	0.310	ug/L	12/18/08
Dibromochloromethane	ND	0.500	0.150	ug/L	12/18/08
1,3-Dichloropropane	ND	0.400	0.120	ug/L	12/18/08
1,2-Dibromoethane	ND	1.00	0.310	ug/L	12/18/08
Carbon tetrachloride	ND	1.00	0.310	ug/L	12/18/08
1,1,1,2-Tetrachloroethane	ND	0.500	0.150	ug/L	12/18/08
Chloroform	ND	1.00	0.300	ug/L	12/18/08
Bromobenzene	ND	1.00	0.310	ug/L	12/18/08
Chloromethane	ND	1.00	0.310	ug/L	12/18/08
1,2,3-Trichloropropane	ND	1.00	0.310	ug/L	12/18/08
Bromomethane	ND	3.00	0.940	ug/L	12/18/08
Bromochloromethane	ND	1.00	0.310	ug/L	12/18/08
Vinyl chloride	ND	1.00	0.310	ug/L	12/18/08
Dichlorodifluoromethane	ND	1.00	0.310	ug/L	12/18/08



SGS Ref.# Client Name 876587 Method Blank

The Environmental Company, Inc. (TEC)

9121-003 Red Hill BFSF Project Name/# Matrix

Water (Surface, Eff., Ground)

**Printed Date/Time** Prep Batch

12/22/2008 11:45 VXX19107

Method Date

SW5030B 12/18/2008

Parameter	Results	Reporting/Control Limit	MDL	Units	Analysis Date
Volatile Gas Chromatography/N	Mass Spectro	oscopy			
Chloroethane	ND	1.00	0.310	ug/L	12/18/08
sec-Butylbenzene	ND	1.00	0.310	ug/L ug/L	12/18/08
Bromodichloromethane	ND	0.500	0.150	ug/L ug/L	12/18/08
1,1-Dichloroethene	ND	1.00	0.130	ug/L	12/18/08
2-Butanone (MEK)	ND	10.0	3.10	ug/L	12/18/08
Methylene chloride	ND	5.00	1.00	ug/L ug/L	12/18/08
Trichlorofluoromethane	ND	1.00	0.310	ug/L ug/L	12/18/08
P & M -Xylene	ND	2.00	0.620	ug/L ug/L	12/18/08
Naphthalene	ND ND	2.00	0.620	ug/L	12/18/08
o-Xylene	ND ND	1.00	0.310	ug/L ug/L	12/18/08
Bromoform	ND	1.00	0.310	ug/L ug/L	12/18/08
1-Chlorohexane	ND ND	1.00	0.310	ug/L ug/L	12/18/08
1,2,4-Trimethylbenzene	ND ND	1.00	0.310	ug/L ug/L	12/18/08
tert-Butylbenzene	ND	1.00	0.310	ug/L ug/L	12/18/08
1,1,1-Trichloroethane	ND ND	1.00	0.310	ug/L ug/L	12/18/08
1,1-Dichloroethane	ND ND	1.00	0.310	ug/L	12/18/08
2-Chlorotoluene	ND ND	1.00	0.310	ug/L ug/L	12/18/08
Trichloroethene	ND ND	1.00	0.310	ug/L	12/18/08
trans-1,2-Dichloroethene	ND ND	1.00	0.310	ug/L	12/18/08
1,2-Dichlorobenzene	ND ND	1.00	0.310	ug/L ug/L	12/18/08
2,2-Dichloropropane	ND ND	1.00	0.310	ug/L	12/18/08
Hexachlorobutadiene	ND ND	1.00	0.310		12/18/08
	ND ND	1.00	0.310	ug/L ug/L	12/18/08
Isopropylbenzene (Cumene)	ND ND	1.00	0.310	_	12/18/08
1,2-Dichloropropane	ND ND	1.00	0.310	ug/L	12/18/08
1,1-Dichloropropene	ND ND	1.00		ug/L	12/18/08
1,1,2-Trichloroethane		1.00	0.310	ug/L	12/18/08
1,3-Dichlorobenzene	ND	1.00	0.310	ug/L	12/18/08
1,2,3-Trichlorobenzene	ND	1.00	0.310	ug/L	12/16/06
Surrogates					
1,2-Dichloroethane-D4 <surr></surr>	114	73-120		%	12/18/08
Toluene-d8 <surr></surr>	99.4	80-120		%	12/18/08
4-Bromofluorobenzene <surr></surr>	103	76-120		%	12/18/08

VMS10317 Batch Method SW8260B

Instrument HP 5890 Series II MS3 VNA



SGS Ref.# 876588 Lab Control Sample

The Environmental Company, Inc. (TEC)

LCS

LCS

LCSD 32.9

LCSD 30.1

33.3

29.5

111

110

98

100

(80-125)

(80-120)

1

2

(<20)

(<20)

Project Name/# 9121-003 Red Hill BFSF

Prep Batch VXX19107 876589 Lab Control Sample Duplicate Method SW5030BClient Name

12/22/2008

12/18/2008

30 ug/L

30 ug/L

30 ug/L

30 ug/L

12/18/2008

12/18/2008

12/18/2008

12/18/2008

**Printed Date/Time** 

Date

11:45

Matrix Water (Surface, Eff., Ground) QC results affect the following production samples:

 $1086721001,\, 1086721002,\, 1086721003$ 

4-Isopropyltoluene

cis-1,3-Dichloropropene

Parameter		QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chromatograp	hy/Mass Sp	ectrosc	ору					
Benzene	LCS	32.9	110	(80-120)			30 ug/L	12/18/2008
	LCSD	33.1	110		1	(< 20)	30 ug/L	12/18/2008
Toluene	LCS	32.5	108	(77-120)			30 ug/L	12/18/2008
	LCSD	32.2	107		1	(< 20)	30 ug/L	12/18/2008
Ethylbenzene	LCS	32.7	109	(80-120)			30 ug/L	12/18/2008
	LCSD	32.7	109		0	(< 20)	30 ug/L	12/18/2008
n-Butylbenzene	LCS	33.4	111	(80-124)			30 ug/L	12/18/2008
	LCSD	33.1	110		1	(< 20)	30 ug/L	12/18/2008
1,4-Dichlorobenzene	LCS	32.6	109	(80-120)			30 ug/L	12/18/2008
	LCSD	32.2	107		1	(< 20)	30 ug/L	12/18/2008
1,2-Dichloroethane	LCS	32.6	109	(80-129)			30 ug/L	12/18/2008
	LCSD	32.4	108		1	(< 20)	30 ug/L	12/18/2008
1,3,5-Trimethylbenzene	LCS	33.3	111	(80-128)			30 ug/L	12/18/2008
	LCSD	32.8	109		2	(< 20)	30 ug/L	12/18/2008
4-Chlorotoluene	LCS	33.2	111	(79-128)			30 ug/L	12/18/2008
	LCSD	32.9	110		1	(< 20)	30 ug/L	12/18/2008
Chlorobenzene	LCS	30.6	102	(80-120)			30 ug/L	12/18/2008
	LCSD	30.5	102		0	(< 20)	30 ug/L	12/18/2008
4-Methyl-2-pentanone (MIBK)	LCS	93.0	103	( 69-134 )			90 ug/L	12/18/2008
	LCSD	98.2	109		5	(< 20)	90 ug/L	12/18/2008
cis-1,2-Dichloroethene	LCS	34.4	115	(80-125)			30 ug/L	12/18/2008
	LCSD	34.1	114		1	(< 20)	30 ug/L	12/18/2008



Client Name

SGS Ref.# 876588 Lab Control Sample

> 876589 Lab Control Sample Duplicate

The Environmental Company, Inc. (TEC)

Project Name/# 9121-003 Red Hill BFSF

Matrix Water (Surface, Eff., Ground) Printed Date/Time Prep Batch

12/22/2008 VXX19107 11:45

SW5030B

Method Date 12/18/2008

Parameter		QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chromatograp	hy/Mass Sp	ectrosc	ору					
n-Propylbenzene	LCS	33.1	110	(80-129)			30 ug/L	12/18/2008
	LCSD	32.6	109		2	(< 20)	30 ug/L	12/18/2008
Styrene	LCS	30.5	102	(80-120)			30 ug/L	12/18/2008
•	LCSD	30.8	103		1	(< 20)	30 ug/L	12/18/2008
Dibromomethane	LCS	31.4	105	(80-120)			30 ug/L	12/18/2008
	LCSD	32.5	108	, ,	3	(<20)	30 ug/L	12/18/2008
trans-1,3-Dichloropropene	LCS	32.5	108	(80-124)			30 ug/L	12/18/2008
, 1 1	LCSD		110	,	2	(<20)	30 ug/L	12/18/2008
1,2,4-Trichlorobenzene	LCS	30.7	102	(80-120)			30 ug/L	12/18/2008
,,		30.3	101	( )	1	(< 20)	30 ug/L	12/18/2008
Acetone	LCS	94.2	105	(50-135)			90 ug/L	12/18/2008
		101	112	(50 150)	7	(< 20)	90 ug/L	12/18/2008
1,1,2,2-Tetrachloroethane	LCS	32.1	107	(76-123)			30 ug/L	12/18/2008
1,1,2,2		33.0	110	( / 0 120 )	3	(< 20)	30 ug/L	12/18/2008
1,2-Dibromo-3-chloropropane	LCS	34.8	116	(73-130)			30 ug/L	12/18/2008
i, 2 Biotomo 5 Cinotopropune		33.9	113	(75 150)	3	(< 20)	30 ug/L	12/18/2008
Methyl-t-butyl ether	LCS	49.6	110	(80-120)			45 ug/L	12/18/2008
Wethyr-t-outyr ether		49.9	111	(00-120)	1	(< 20)	45 ug/L 45 ug/L	12/18/2008
Tetrachloroethene	LCS	31.2	104	(79-122)			20. ug/I	12/18/2008
renaemoroemene	LCSD		104	(79-122)	0	(< 20)	30 ug/L 30 ug/L	12/18/2008
Dibromochloromothono	LCS	22.0	110	(80-120)			20/I	12/19/2009
Dibromochloromethane	LCSD	33.0 33.5	110	( 80-120 )	2	(< 20)	30 ug/L 30 ug/L	12/18/2008 12/18/2008
12 D: 11				(00.121)			20 /7	12/10/2000
1,3-Dichloropropane	LCS LCSD	33.1 33.4	110 111	(80-121)	1	(< 20)	30 ug/L 30 ug/L	12/18/2008 12/18/2008
10.07				(00.100)		, ,		
1,2-Dibromoethane	LCS LCSD	32.5 32.9	108 110	(80-120)	1	(< 20)	30 ug/L 30 ug/L	12/18/2008 12/18/2008
						( /		
Carbon tetrachloride	LCS LCSD	32.9 32.9	110 110	(80-126)	0	(< 20)	30 ug/L 30 ug/L	12/18/2008 12/18/2008
	LCSD	34.9	110		U	( > 20 )	50 ug/L	12/10/2000



SGS Ref.# 876588 Lab Control Sample

876589 Lab Control Sample Duplicate

Client Name The Environmental Company, Inc. (TEC)
Project Name/# 9121-003 Red Hill BFSF

Matrix Water (Surface, Eff., Ground)

 Printed Date/Time
 12/22/2008
 11:45

 Prep
 Batch
 VXX19107

 Method
 SW5030B

Date 12/18/2008

Parameter		QC Results	Pct Recov	LCS/LCSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chromatogra	aphy/Mass Sr	pectrosc	ору					
1,1,1,2-Tetrachloroethane	LCS	32.4	108	(80-120)			30 ug/L	12/18/2008
	LCSD	32.6	109		1	(< 20)	30 ug/L	12/18/2008
Chloroform	LCS	32.2	107	(80-124)			30 ug/L	12/18/2008
	LCSD	33.0	110		3	(< 20)	30 ug/L	12/18/2008
Bromobenzene	LCS	31.5	105	(80-120)			30 ug/L	12/18/2008
	LCSD	31.4	105		0	(< 20)	30 ug/L	12/18/2008
Chloromethane	LCS	32.7	109	(67-125)			30 ug/L	12/18/2008
	LCSD		106	,	3	(<20)	30 ug/L	12/18/2008
1,2,3-Trichloropropane	LCS	31.9	106	(80-120)			30 ug/L	12/18/2008
1 1	LCSD	32.9	110	,	3	(< 20)	30 ug/L	12/18/2008
Bromomethane	LCS	32.2	107	(30-140)			30 ug/L	12/18/2008
	LCSD		105	,	2	(< 20)	30 ug/L	12/18/2008
Bromochloromethane	LCS	32.2	107	(77-129)			30 ug/L	12/18/2008
	LCSD		108	,	1	(<20)	30 ug/L	12/18/2008
Vinyl chloride	LCS	32.9	110	(72-145)			30 ug/L	12/18/2008
,	LCSD		107	,	3	(< 20)	30 ug/L	12/18/2008
Dichlorodifluoromethane	LCS	32.6	109	(62-153)			30 ug/L	12/18/2008
	LCSD		106	,	3	(< 20)	30 ug/L	12/18/2008
Chloroethane	LCS	34.1	114	(67-133)			30 ug/L	12/18/2008
	LCSD		112	,	2	(< 20)	30 ug/L	12/18/2008
sec-Butylbenzene	LCS	33.6	112	(80-120)			30 ug/L	12/18/2008
,	LCSD		111	,	1	(< 20)	30 ug/L	12/18/2008
Bromodichloromethane	LCS	31.1	104	(80-120)			30 ug/L	12/18/2008
	LCSD		105		1	(<20)	30 ug/L	12/18/2008
1,1-Dichloroethene	LCS	34.2	114	(76-130)			30 ug/L	12/18/2008
,	LCSD	34.4	115	( )	1	(< 20)	30 ug/L	12/18/2008
2-Butanone (MEK)	LCS	93.1	103	(66-136)			90 ug/L	12/18/2008
• /				` ,			<i>G</i> -	



Client Name

876588 Lab Control Sample

876589 Lab Control Sample Duplicate

The Environmental Company, Inc. (TEC)

Project Name/#

9121-003 Red Hill BFSF

Water (Surface, Eff., Ground)

**Printed Date/Time** Prep Batch Method

Date

12/22/2008 VXX19107 11:45

SW5030B 12/18/2008

Matrix RPD Pct LCS/LCSD Spiked Analysis RPD Parameter Results Limits Limits Recov Amount Date Volatile Gas Chromatography/Mass Spectroscopy 9 LCSD 102 (< 20)90 ug/L 12/18/2008 114 Methylene chloride 110 LCS 33.1 (63-131)30 ug/L 12/18/2008 5 (< 20)30 ug/L LCSD 31.4 105 12/18/2008 Trichlorofluoromethane LCS 35.4 118 (68-145)30 ug/L12/18/2008 3 (< 20)30 ug/L 12/18/2008 LCSD 34.2 114 P & M -Xylene LCS 65.5 109 (80-120)60 ug/L 12/18/2008 2 LCSD 66.5 111 (< 20)60 ug/L 12/18/2008 Naphthalene LCS 30.8 103 (75-120)30 ug/L 12/18/2008 99 3 (< 20)30 ug/L 12/18/2008 LCSD 29.8 LCS o-Xylene 32.9 110 (80-120)30 ug/L 12/18/2008 1 (< 20)30 ug/L 12/18/2008 LCSD 32.7 109 Bromoform LCS 34.0 113 (80-120)30 ug/L 12/18/2008 (< 20)LCSD 34.4 115 1 30 ug/L 12/18/2008 1-Chlorohexane LCS 50.8 113 (70-125)45 ug/L 12/18/2008 1 (< 20)45 ug/L 12/18/2008 LCSD 50.4 112 1,2,4-Trimethylbenzene 109 (80-125)LCS 32.8 30 ug/L12/18/2008 LCSD 32.4 108 1 (< 20)30 ug/L 12/18/2008 tert-Butylbenzene LCS 34.5 115 (80-122)30 ug/L 12/18/2008 30 ug/L LCSD 34.1 114 1 (< 20)12/18/2008 1,1,1-Trichloroethane LCS 31.0 103 (80-122)30 ug/L12/18/2008 1 (< 20)30 ug/L 12/18/2008 LCSD 31.3 104 1,1-Dichloroethane LCS 33.3 111 (80-120)30 ug/L 12/18/2008 1 30 ug/L LCSD 32.9 110 (< 20)12/18/2008 2-Chlorotoluene LCS 32.5 108 (80-125)30 ug/L12/18/2008 1 (< 20)30 ug/L 12/18/2008 LCSD 32.3 108 Trichloroethene 108 LCS 32.4 (80-125)30 ug/L 12/18/2008 LCSD 32.9 110 2 (< 20)30 ug/L 12/18/2008



4-Bromofluorobenzene <surr>

LCS

LCSD

SGS Ref.#

876588 Lab Control Sample

876589 Lab Control Sample Duplicate

The Environmental Company, Inc. (TEC)

Client Name Project Name/#

9121-003 Red Hill BFSF

Matrix Water (Surface, Eff., Ground)

Printed Date/Time Prep Batch 12/22/2008 VXX19107 11:45

12/18/2008 12/18/2008

Method SW5030B

Date 12/18/2008

LCS/LCSD RPD Pct Spiked Analysis RPD Parameter Results Limits Limits Recov Date Amount Volatile Gas Chromatography/Mass Spectroscopy trans-1,2-Dichloroethene LCS 32.4 108 (79-132)30 ug/L 12/18/2008 LCSD 32.2 107 1 30 ug/L 12/18/2008 (< 20)1,2-Dichlorobenzene LCS 31.6 105 (80-120)30 ug/L 12/18/2008 LCSD 31.6 105 0 (< 20)30 ug/L 12/18/2008 LCS 34.5 115 2,2-Dichloropropane (80-132)30 ug/L 12/18/2008 LCSD 33.8 113 2 (< 20)30 ug/L 12/18/2008 LCS 32.2 107 Hexachlorobutadiene (77-125)30 ug/L 12/18/2008 LCSD 31.7 106 1 (< 20)30 ug/L 12/18/2008 Isopropylbenzene (Cumene) LCS 34.3 114 (80-121)30 ug/L 12/18/2008 1 (< 20)30 ug/L 12/18/2008 LCSD 34.1 114 107 1,2-Dichloropropane LCS 32.0 (80-121)30 ug/L 12/18/2008 2 LCSD 32.7 109 (< 20)30 ug/L 12/18/2008 LCS 32.4 108 (80-122)1,1-Dichloropropene 30 ug/L 12/18/2008 108 0 (< 20)30 ug/L 12/18/2008 LCSD 32.3 LCS 108 1,1,2-Trichloroethane 32.3 (77-120)30 ug/L 12/18/2008 LCSD 32.5 108 0 (< 20)30 ug/L 12/18/2008 107 1,3-Dichlorobenzene LCS 32.1 (80-120)30 ug/L 12/18/2008 0 12/18/2008 107 (< 20)30 ug/L LCSD 32.0 1,2,3-Trichlorobenzene LCS 31.0 103 (77-120)30 ug/L 12/18/2008 LCSD 30.6 102 1 (< 20)30 ug/L 12/18/2008 Surrogates 1,2-Dichloroethane-D4 <surr> LCS 99 (73-120)12/18/2008 2 12/18/2008 **LCSD** 100 Toluene-d8 <surr> 99 LCS (80-120)12/18/2008 **LCSD** 100 1 12/18/2008

98

98

(76-120)

0



 SGS Ref.#
 876588
 Lab Control Sample
 Printed Date/Time
 12/22/2008
 11:45

876589 Lab Control Sample Duplicate Prep Batch VXX19107
Client Name The Environmental Company, Inc. (TEC) Method SW5030B

Project Name/# 9121-003 Red Hill BFSF Date 12/18/2008

Matrix Water (Surface, Eff., Ground)

Parameter QC Pct LCS/LCSD RPD Spiked Analysis
Parameter Results Recov Limits RPD Limits Amount Date

#### Volatile Gas Chromatography/Mass Spectroscopy

Batch VMS10317 Method SW8260B

Instrument HP 5890 Series II MS3 VNA



876591 876592 Matrix Spike

Matrix Spike Duplicate

Printed Date/Time

12/22/2008 11:45

Prep Batch VXX19107 Method Volatiles Extraction AFCEE 3.1

Date

12/18/2008

Original 876590

Matrix Water (Surface, Eff., Ground)

QC results affect the following production samples: 1086721001, 1086721002, 1086721003

Parameter	Qualifiers	Original Result	QC Result	Pct Recov	MS/MSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chron	matography/I	Mass Spe	ctroscopy						
Benzene	MS	ND	33	110	(80-120)			30.0	ug/L 12/18/2008
	MSD		34.2	114		4	(< 20)	30.0	ug/L 12/18/2008
Toluene	MS	ND	32.2	107	(77-120)			30.0	ug/L 12/18/2008
	MSD		33.4	111		4	(< 20)	30.0	ug/L 12/18/2008
Ethylbenzene	MS	ND	33.1	110	(80-120)			30.0	ug/L 12/18/2008
	MSD		34.1	114		3	(< 20)	30.0	ug/L 12/18/2008
n-Butylbenzene	MS	ND	35.1	117	(80-124)			30.0	ug/L 12/18/2008
	MSD		37.7	126*		7	(< 20)	30.0	ug/L 12/18/2008
1,4-Dichlorobenzene	MS	ND	33	110	(80-120)			30.0	ug/L 12/18/2008
	MSD		35.3	118		7	(< 20)	30.0	ug/L 12/18/2008
1,2-Dichloroethane	MS	ND	38	127	(80-129)			30.0	ug/L 12/18/2008
	MSD		34.5	115		10	(< 20)	30.0	ug/L 12/18/2008
1,3,5-Trimethylbenzene	MS	ND	33.8	113	(80-128)			30.0	ug/L 12/18/2008
	MSD		35.4	118		5	(< 20)	30.0	ug/L 12/18/2008
4-Chlorotoluene	MS	ND	33.7	112	(79-128)			30.0	ug/L 12/18/2008
	MSD		35.3	118		5	(< 20)	30.0	ug/L 12/18/2008
Chlorobenzene	MS	ND	30.5	102	(80-120)			30.0	ug/L 12/18/2008
	MSD		31.5	105		3	(< 20)	30.0	ug/L 12/18/2008
4-Methyl-2-pentanone (M	IIBK) MS	ND	102	113	(69-134)			90.0	ug/L 12/18/2008
	MSD		103	115		1	(< 20)	90.0	ug/L 12/18/2008
cis-1,2-Dichloroethene	MS	ND	38.6	129*	(80-125)			30.0	ug/L 12/18/2008
	MSD		36.3	121		6	(< 20)	30.0	ug/L 12/18/2008
4-Isopropyltoluene	MS	ND	34.1	114	(80-125)			30.0	ug/L 12/18/2008
	MSD		36.4	121		7	(< 20)	30.0	ug/L 12/18/2008
cis-1,3-Dichloropropene	MS	ND	31.7	106	(80-120)			30.0	ug/L 12/18/2008
	MSD		31.9	106		1	(< 20)	30.0	ug/L 12/18/2008
n-Propylbenzene	MS	ND	33.8	113	(80-129)			30.0	ug/L 12/18/2008
	MSD		35.4	118		5	(< 20)	30.0	ug/L 12/18/2008
Styrene	MS	ND	31.2	104	(80-120)			30.0	ug/L 12/18/2008
	MSD		31.5	105		1	(< 20)	30.0	ug/L 12/18/2008
Dibromomethane	MS	ND	37.2	124*	(80-120)			30.0	ug/L 12/18/2008
	MSD		34.1	114		9	(< 20)	30.0	ug/L 12/18/2008
trans-1,3-Dichloropropen	e MS	ND	33.6	112	(80-124)			30.0	ug/L 12/18/2008
	MSD		33.5	112		0	(< 20)	30.0	ug/L 12/18/2008
1,2,4-Trichlorobenzene	MS	ND	31.7	106	(80-120)			30.0	ug/L 12/18/2008
	MSD		34.1	114		7	(< 20)	30.0	ug/L 12/18/2008
Acetone	MS	ND	104	116	(50-135)			90.0	ug/L 12/18/2008



876591 876592 Matrix Spike

Matrix Spike Duplicate

Printed Date/Time
Prep Batch

e/Time 12/22/2008 11:45 Batch VXX19107

Method Volatiles Extraction AFCEE 3.1

Date 12/18/2008

Original 876590

Matrix Water (Surface, Eff., Ground)

Parameter Qua	Original lifiers Result	QC Result	Pct MS/MSD Recov Limits	RPD RPD Limits	Spiked Analysis Amount Date
Volatile Gas Chromat	ography/Mass Sm	pectroscopy			
	MSD	102	114	2 (< 20)	90.0 ug/L 12/18/2008
1,1,2,2-Tetrachloroethane	MS ND	34.1	114 (76-123)		30.0 ug/L 12/18/2008
	MSD	34.8	116	2 (< 20)	30.0 ug/L 12/18/2008
1,2-Dibromo-3-chloropropan	e MS ND	35.2	117 (73-130)		30.0 ug/L 12/18/2008
	MSD	36.1	120	2 (< 20)	30.0 ug/L 12/18/2008
Methyl-t-butyl ether	MS ND	55	122* (80-120)		45.0 ug/L 12/18/2008
	MSD	53.2	118	3 (< 20)	45.0 ug/L 12/18/2008
Tetrachloroethene	MS ND	32.4	108 (79-122)		30.0 ug/L 12/18/2008
	MSD	32.9	110	2 (< 20)	30.0 ug/L 12/18/2008
Dibromochloromethane	MS ND	34.5	115 (80-120)		30.0 ug/L 12/18/2008
	MSD	33.4	111	3 (< 20)	30.0 ug/L 12/18/2008
1,3-Dichloropropane	MS ND	35.1	117 (80-121)		30.0 ug/L 12/18/2008
	MSD	34.4	115	2 (< 20)	30.0 ug/L 12/18/2008
1,2-Dibromoethane	MS ND	34.5	115 (80-120)		30.0 ug/L 12/18/2008
	MSD	33.8	113	2 (< 20)	30.0 ug/L 12/18/2008
Carbon tetrachloride	MS ND	36.5	122 (80-126)		30.0 ug/L 12/18/2008
	MSD	35.1	117	4 (< 20)	30.0 ug/L 12/18/2008
1,1,1,2-Tetrachloroethane	MS ND	33.1	110 (80-120)		30.0 ug/L 12/18/2008
	MSD	32.6	109	2 (< 20)	30.0 ug/L 12/18/2008
Chloroform	MS ND	38.5	128* (80-124)		30.0 ug/L 12/18/2008
	MSD	35.1	117	9 (< 20)	30.0 ug/L 12/18/2008
Bromobenzene	MS ND	32.4	108 (80-120)		30.0 ug/L 12/18/2008
	MSD	33.6	112	4 (< 20)	30.0 ug/L 12/18/2008
Chloromethane	MS ND	28.8	96 (67-125)		30.0 ug/L 12/18/2008
	MSD	31.3	104	8 (< 20)	30.0 ug/L 12/18/2008
1,2,3-Trichloropropane	MS ND	32.7	109 (80-120)		30.0 ug/L 12/18/2008
• •	MSD	34.8	116	6 (< 20)	30.0 ug/L 12/18/2008
Bromomethane	MS ND	32	107 (30-140)		30.0 ug/L 12/18/2008
	MSD	33.8	113	6 (< 20)	30.0 ug/L 12/18/2008
Bromochloromethane	MS ND	34	113 (77-129)		30.0 ug/L 12/18/2008
	MSD	33.8	113	1 (< 20)	30.0 ug/L 12/18/2008
Vinyl chloride	MS ND	31	103 (72-145)		30.0 ug/L 12/18/2008
•	MSD	32.0	107	3 (< 20)	30.0 ug/L 12/18/2008
Dichlorodifluoromethane	MS ND	33.3	111 (62-153)		30.0 ug/L 12/18/2008
	MSD	33.8	113	2 (< 20)	30.0 ug/L 12/18/2008
Chloroethane	MS ND	49.8	166* (67-133)		30.0 ug/L 12/18/2008
	MSD	51.3	171*	3 (< 20)	30.0 ug/L 12/18/2008
sec-Butylbenzene	MS ND	34.1	114 (80-120)		30.0 ug/L 12/18/2008
•	MSD	35.9	120	5 (< 20)	30.0 ug/L 12/18/2008
Bromodichloromethane	MS ND	34.5	115 (80-120)		30.0 ug/L 12/18/2008
2. Omoundmondmondmondmon	1110 1110	51.5	115 (00 120)		J0.0 48/L 12/10/2000



876591 876592 Matrix Spike

Matrix Spike Duplicate

Printed Date/Time Prep

12/22/2008 11:45 Batch

VXX19107

Method

Volatiles Extraction AFCEE 3.1

Date 12/18/2008

Original 876590

Matrix Water (Surface, Eff., Ground)

Parameter	Qualifiers	Original Result	QC Result	Pct Recov	MS/MSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chr	omatography/l	Mass Spe	ectroscopy						
	MSD		32.6	109		5	(< 20)	30.0 u	g/L 12/18/2008
1,1-Dichloroethene	MS	ND	37.4	125	(76-130)		,		g/L 12/18/2008
,	MSD		35.8	119	· · · · · ·	5	(< 20)		g/L 12/18/2008
2-Butanone (MEK)	MS	ND	105	117	(66-136)				g/L 12/18/2008
, ,	MSD		105	116		0	(< 20)		g/L 12/18/2008
Methylene chloride	MS	ND	31.4	105	(63-131)			30.0 u	g/L 12/18/2008
	MSD		32.3	108		3	(< 20)		g/L 12/18/2008
Trichlorofluoromethane	MS	ND	38	127	(68-145)			30.0 u	g/L 12/18/2008
	MSD		37.4	125		2	(< 20)		g/L 12/18/2008
P & M -Xylene	MS	ND	66.4	111	(80-120)			60.0 u	g/L 12/18/2008
	MSD		68.6	114		3	(< 20)	60.0 u	g/L 12/18/2008
Naphthalene	MS	ND	31.3	104	(75-120)			30.0 u	g/L 12/18/2008
	MSD		34.1	114		8	(< 20)	30.0 u	g/L 12/18/2008
o-Xylene	MS	ND	32.9	110	(80-120)			30.0 u	g/L 12/18/2008
	MSD		33.9	113		3	(< 20)	30.0 u	g/L 12/18/2008
Bromoform	MS	ND	34.6	115	(80-120)			30.0 u	g/L 12/18/2008
	MSD		34.2	114		1	(< 20)	30.0 u	g/L 12/18/2008
1-Chlorohexane	MS	ND	53.6	119	(70-125)			45.0 u	g/L 12/18/2008
	MSD		55.6	123		4	(< 20)	45.0 u	g/L 12/18/2008
1,2,4-Trimethylbenzene	MS	ND	33.2	111	(80-125)			30.0 u	g/L 12/18/2008
	MSD		34.8	116		5	(< 20)	30.0 u	g/L 12/18/2008
tert-Butylbenzene	MS	ND	34.8	116	(80-122)			30.0 u	g/L 12/18/2008
	MSD		36.8	123*		6	(< 20)	30.0 u	g/L 12/18/2008
1,1,1-Trichloroethane	MS	ND	34.5	115	(80-122)			30.0 u	g/L 12/18/2008
	MSD		33.0	110		5	(< 20)	30.0 u	g/L 12/18/2008
1,1-Dichloroethane	MS	ND	37	123*	(80-120)			30.0 u	g/L 12/18/2008
	MSD		34.8	116		6	(< 20)	30.0 u	g/L 12/18/2008
2-Chlorotoluene	MS	ND	32.8	109	(80-125)			30.0 u	g/L 12/18/2008
	MSD		34.8	116		6	(< 20)		g/L 12/18/2008
Trichloroethene	MS	ND	33	110	(80-125)			30.0 u	g/L 12/18/2008
	MSD		34.4	115		4	(< 20)	30.0 u	g/L 12/18/2008
trans-1,2-Dichloroethen		ND	38.4	128	(79-132)			30.0 u	g/L 12/18/2008
	MSD		34.2	114		12	(< 20)	30.0 u	g/L 12/18/2008
1,2-Dichlorobenzene	MS	ND	32.4	108	(80-120)				g/L 12/18/2008
	MSD		33.9	113		4	(< 20)		g/L 12/18/2008
2,2-Dichloropropane	MS	ND	39.7	132	(80-132)				g/L 12/18/2008
	MSD		38.6	129		3	(< 20)		g/L 12/18/2008
Hexachlorobutadiene	MS	ND	33.8	113	(77-125)				g/L 12/18/2008
	MSD		35.9	120		6	(< 20)		g/L 12/18/2008
Isopropylbenzene (Cum	nene) MS	ND	34.8	116	(80-121)			30.0 u	g/L 12/18/2008



876591

Matrix Spike

876592 Matrix Spike Duplicate

Printed Date/Time

12/22/2008 11:45

Prep Batch

VXX19107

Method

Volatiles Extraction AFCEE 3.1

Date

12/18/2008

Original

876590

Matrix Water (Surface, Eff., Ground)

Parameter	Qualifiers	Original Result	QC Result	Pct Recov	MS/MSD Limits	RPD	RPD Limits	Spiked Amount	Analysis Date
Volatile Gas Chro	matography/	Mass Spe	ectroscopy						
	MSE	)	35.5	118		2	(< 20)	30.0	ug/L 12/18/2008
1,2-Dichloropropane	MS	ND	33.3	111	(80-121)			30.0	ug/L 12/18/2008
	MSE	)	34.3	114		3	(< 20)	30.0	ug/L 12/18/2008
1,1-Dichloropropene	MS	ND	34.4	115	(80-122)			30.0	ug/L 12/18/2008
	MSE	)	34.8	116		1	(< 20)	30.0	ug/L 12/18/2008
1,1,2-Trichloroethane	MS	ND	33.7	112	(77-120)			30.0	ug/L 12/18/2008
	MSE	)	33.4	111		1	(< 20)	30.0	ug/L 12/18/2008
1,3-Dichlorobenzene	MS	ND	32.7	109	(80-120)			30.0	ug/L 12/18/2008
	MSE	)	34.6	115		6	(< 20)	30.0	ug/L 12/18/2008
1,2,3-Trichlorobenzene	MS	ND	31.2	104	(77-120)			30.0	ug/L 12/18/2008
	MSE	)	33.6	112		7	(< 20)	30.0	ug/L 12/18/2008
Surrogates									
1,2-Dichloroethane-D4 <	surr> MS		34.2	114	(73-120)				12/18/2008
	MSE	)	29.9	100		13			12/18/2008
Toluene-d8 <surr></surr>	MS		29.7	99	(80-120)				12/18/2008
	MSE	)	29.3	98		2			12/18/2008
4-Bromofluorobenzene <	surr> MS		29.3	98	(76-120)				12/18/2008
	MSE	)	29.6	99		1			12/18/2008

Batch VMS10317 Method SW8260B

Instrument HP 5890 Series II MS3 VNA



## CHAIN OF CUSTODY RECORD SGS Environmental Services Inc.



#### Locations Nationwide

Macka

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Maryland

Louisiana

North Carolina

West Virginia

www.us.sgs.com

CLIENT:	TEC INC.				SGS Ref	ference #:													,
CONTACT:	Jeff Hart PHO	ONE NO:	303.2 <b>7</b> 3.0231		<u>L</u>											page		of	-
PROJECT:	9121-003 sitt	E/PWSID#:	Red Hill BFSF			Preserv. Used	/ <sub>X</sub> O		\s\cdot\)	$\angle$	, HILO	//			$\angle$				
REPORTS TO:	Jeff Hart ema		ecinc.com man@tecinc.	com	# C O	SAMPLE TYPE C =	e	<u>6</u>											
INVOICE TO:		OTE #:			T A I N	COMP G =	тРн-GRO (8015В)	TPH-DRO (8015B)	VOC's (8260B)	PAH's (8270-SIM)	Pb (6020)								
LAB NO.	SAMPLE IDENTIFICATION	DATE	TIME	MATRIX	E R S	GRAB	5-нат	ТРН-D	NOC's	PAH's	Diss P							REMARKS	
0 A·c 2   3	RHMW2254-WG13B	12/16/08	1005	Water	3				Х					<u> </u>			_		
	RHMWA01-WG13B	12/11/08	1205	Water	3	ļ			Х				_				<u> </u>		
(3) V	TB01-WG13B	12/16/08	0805	Water	3		ļ		Х					-	ļi	·	- -		
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Collected/Relinquis	DU LEE	Date 12/16/08	1300	Received By:						ng Carri						Samples G Temperatu	Received ( ††) : ure °C:	Cold? YES NO	<b>,</b> T
Relinquished By:	(2)	Date	Time	Received By:					Special	Delive	rable R	equiremen	s:			Chain of	Custody BROKE	Seal: (Circle)	
Relinquished By:	(3)	Date	Time	Received By:					Reques	sted Tu	rnaround	l Time ar	d-or Spe	cial Instr	uctions:				
Relinquished By:	(4)	Date 12/1/08	Time US 5	Received For Lab	oratory By:	2													
3180 Pege	tter Drive Anchorage, AK 99518 Tel: (907) r Road Fairbanks, AK 99701 Tel: (907) 47 Island Access Rd., Unit 1B Honolulu, HI 961	562-2343 Fax: (90° 4-8656 Fax: (907)	74-9685	15-2287			Greent	orier Str	reet Ch	arlestor	n, WV :	25311 Tel	(304) 3	346-0725	Fax:	463-330 (304) 346 10) 350-15	-0761		

SGS

# 1086721

#### SAMPLE RECEIPT FORM

SGS WO#:

Yes	No NA	A	4.
	V	Are samples RUSH, priority or w/in 72 hrs of hold time?	TAT (circle one): Standard -or- Rush
	— <del>-</del>	If yes, have you done <i>e-mail ALERT notification</i> ?	Received Date: 12.17-08
	<u> </u>	Are samples within 24 hrs. of hold time or due date?	Received Time: //55
	ーし	If yes, have you also <i>spoken with</i> supervisor?	Is date/time conversion necessary?
		Archiving bottles (if req'd): Are they properly marked?	# of hours to AK Local Time:
		Are there any <b>problems</b> ? PM Notified?	Thermometer ID: 696
		Were samples preserved correctly and pH-verified?	Cooler ID Temp Blank Cooler Temp
1			/ 2.3 °C 3./ °C
			°C °C
,			°C
	<i></i>	If this is for PWS, provide PWSID.	°C
	v _	_ Will courier charges apply?	°C
		Method of payment?	Note: Temperature readings include thermometer correction factors
$\underline{\nu}$ _		_ Data package required? (Level: 1 / 2 / 3 / 4)	Delivery method (circle all that apply): Client /
		Notes:	Alert Courier / UPS / FedEx / USPS / DHL /
		Is this a DoD project? (USACE, Navy, AFCEE)	AA Goldstreak / NAC / ERA / PenAir / Carlile/
XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX	XXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXXX		Lynden / SGS / Other:
		n must be filled out for DoD projects (USACE, Navy, AFCEE)	Airbill # 796/-909/-7035
Yes	No	Is received temperature $4 \pm 2^{\circ}$ C?	Additional Sample Remarks: $(\sqrt{if applicable})$ Extra Sample Volume?
		Exceptions: Samples/Analyses Affected:	Limited Sample Volume?
			MeOH field preserved for volatiles
**************************************			Field-filtered for dissolved
XXXXXXXX XXXXXXX XXXXXXXX			Lab-filtered for dissolved
********* *******	NL	If temperature(s) <0 °C, were containers ice-free? N/A	Ref Lab required?
سا ا		Notify PM immediately of any ice in samples.  Was there an airbill? (Note # above in the right hand column)	Foreign Soil?
7_		Was cooler sealed with custody seals?	
	N	#/ where: 2 FRONT PLACE TOP LIN	This section must be filled if problems are found.
ب		Were seal(s) intact upon arrival?	Yes No
U		Was there a COC with cooler?	Was client notified of problems?
\_V		Was COC sealed in plastic bag & taped inside lid of cooler?	Individual contacted:
1		Was the COC filled out properly?	Via: Phone / Fax / Email (circle one)
	س	Did the COC indicate USACE / Navy / AFCEE project? Did the COC and samples correspond?	Date/Time:
1 - 6		Were all sample packed to prevent breakage?	Reason for contact:
		Packing material: BUMPIF UNAP FULM BLOCK	
<u>ب</u> ا		Were all samples unbroken and clearly labeled?	
X X X X X X X X X X X X X X X X X X X	_ب	Were all samples sealed in separate plastic bags?	
<u></u>	X	Were all VOCs free of headspace and/or MeOH preserved?	
<u> </u>		Were correct container / sample sizes submitted?	Change Order Required?
-		Is sample condition good?  Was copy of CoC, SRF, and custody seals given to PM to fax?	SGS Contact:
		was copy of coc, sre, and custody scals given to rive to tax:	
Notes:			
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## SAMPLE RECEIPT FORM (page 2)



SGS WO#:

						Container Volume							Container Type									Preservative								
#	Container ID	Matrix	Test	OC	TB	11	500 mL	250 mL	125 mL	60 mL				Other	AG	ĐO	HDPE	Nalgene		Coli	Septa	Other	None	HCI	HNO <sub>3</sub>				NaOH	Other
1,2	A.C	1	VOC								6				V						0			٦						
3	A-C	1	VOC	-	レ						3				4						_			<u>_</u>						
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Bottle Totals						9				
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Completed by:

\_\_\_\_I

Date: 17-17-08

From: Origin ID: HIKA (808) 528-1445 BILL WHITMAN TEC INC. 1001 BISHOP STREET, ASB TOWER SUITE 1400 HONOLULU, HI 96813



JCLS111208/20/23

SHIP TO: (907) 562-2343

**BILL THIRD PARTY** 

SAMPLE RECEIVING
SGS Environmental Services
200 W POTTER DR

**ANCHORAGE, AK 99518** 



Ship Date: 16DEC08 ActWgt: 10.0 LB CAD: 1774997/INET8091 Account#: S \*\*\*\*\*\*\*\*\*\*\*

Delivery Address Bar Code



Ref # P# 66: Invoice # PO # Dept #



TRK# 0201

7961 9097 7035

WED - 17DEC AM
PRIORITY OVERNIGHT

**WU CYMA** 

99518 AK-US ANC



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- 2. Fold the printed page along the horizontal line.
- 3. Place label in shipping pouch and affix it to your shipment so that the barcode portion of the label can be read and scanned.

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