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4:11 pm, May 09, 2011 Alameda County Environmental Health

5 May 2011 Project No. 730482302

Ms. Barbara Jakub, P.G. Hazardous Materials Specialist Alameda County Health Care Services Agency 1131 Harbor Bay Parkway, Suite 250 Alameda, California 94502

Subject:

Summary of Soil and Groundwater Analytical Results

Fuel Leak Case Number RO0000201

Hydraulic Electro Service

5812 Hollis Street Emeryville, California

Dear Ms. Jakub:

As a legally authorized representative of Wareham Property Group, and on behalf of Wareham Property Group, I declare, under penalty of perjury, that the information and/or recommendations contained in the attached document titled *Summary of Soil and Groundwater Analytical Results, Fuel Leak Case Number R00000201, Hydraulic Electro Service, 5812 Hollis Street, Emeryville, California,* are true and correct to the best of my knowledge.

Sincerely yours,

Geoffrey B. Sears

WAREHAM PROPERTY GROUP

Attachment



MEMORANDUM

TO:

Ms. Barbara J. Jakub, P. G. - Alameda County Environmental Health

GROVER BUHR

No. 5596

CC:

Mr. Geoff Sears – Wareham

FROM:

Peter J. Cusack, REA

Philip G. Smith, REA

Grover Buhr, P.G.

Treadwell & Rollo, I

DATE:

5 May 2011

PROJECT:

EmeryStation Greenway

5812 Hollis Street Emeryville, California

730482302

SUBJECT:

Summary of Soil and Groundwater Analytical Results

Fuel Leak Case Number RO0000201

Hydraulic Electro Service 5812 Hollis Street

Emeryville, California

Introduction

On behalf of EmeryStation Triangle II LLC, this memorandum presents the analytical results of verification sampling performed at the EmeryStation Greenway project at 5812 Hollis Street (Site) in Emeryville, California. Per our telephone calls and emails, we are transmitting this memorandum with the analytical results of the verification sampling performed at the Site to supplement our 17 September 2010 case closure request and obtain case closure. The Site is bound by Hollis Street to the west, a commercial building to the north (at 5850 Hollis Street), and a public greenway approximately 25 feet wide to the east. The southern tip of the Site is at the intersection of Hollis Street and Powell Street.

The Site is an approximate one-acre triangular-shaped parcel located at the intersection of Hollis and Powell Streets (Figure 1), and was historically operated as a dismantling yard. An underground fuel storage tank (UST) was installed in 1977 in the northern portion of the property for fleet fueling operations. This UST was removed in 1989. Prior to the excavation activities now in progress, the Site was occupied by a paved yard and the foundations and floor slab from a former one-story building that had previously been located in the southern portion of the property.



Excavation Activities

Currently, a four-story commercial building overlying one level of subgrade parking is being constructed at the Site. As part of the Site development plans, the Site has been excavated to approximately 15 feet below ground surface (bgs), with the excavation extending to just beyond the property lines to the east and west. The building will encompass approximately two-thirds of the property. The remaining one-third of the property (the southern portion) will be transferred to the City of Emeryville to tie in to an existing greenway running along the adjacent eastern property boundary.

Numerous investigation activities have been conducted at the Site, and contaminants found at the Site from past use activities included metals (lead) above background levels, some semi-volatile and volatile organic compounds, and petroleum hydrocarbons in the soil and groundwater. Metals and volatile organic compound contamination were found primarily in shallow soil in specific localized areas within the former dismantling yard. Petroleum hydrocarbon-impacted soil and groundwater were found in the immediate vicinity of the former UST. The California Department of Toxic Substances Control (DTSC) is the regulatory oversight agency for the impacted soil not associated with the USTs. Alameda County Environmental Health (ACEH) is the lead oversight agency for the former UST closures, including petroleum hydrocarbon constituents in soil or groundwater associated with the former USTs.

To address Site contamination issues, Treadwell & Rollo prepared a Site Management Plan (SMP) dated 16 February 2010, which describes measures to be taken during redevelopment to ensure that historic use-related substances are removed in a safe and protective manner during construction. This plan was submitted to and approved by DTSC on 21 March 2010; DTSC stated that the proposed scope of work was acceptable and that construction could begin. DTSC also requested a post-excavation report documenting Site soil excavation and disposal, and describing residual soil conditions. Based on a telephone conversation on 11 October 2011 with Ms. Nina Bacey of DTSC regarding sidewall sampling, Ms. Bacey agreed to change the sidewall sampling frequency from collecting a sidewall sample every 25 horizontal feet and 3 vertical feet to a sidewall sample every 50 horizontal feet and vertical 5 feet vertical.

During demolition of the building slab in the southern part of the Site, a second UST was uncovered underneath the northern portion of the slab, in the east-central portion of the Site (Figure 2). The 1,100-gallon UST was removed by Pacific States Environmental Contractors. Inc. of Dublin, California (PSEC) on 20 August 2010 with oversight by ACEH, and approximately 30 cubic yards of soil were excavated to remove petroleum contaminated soil.

On 17 September 2010, a *Case Closure Summary, Hydraulic Electro Service, 5812 Hollis Street, Emeryville, California Fuel Leak Case Number RO0000201* was prepared by Treadwell & Rollo and submitted to ACEH requesting case closure of the Site. In December 2010, a meeting was held at ACEH to discuss the Site redevelopment and what was needed to obtain case closure of the former leaking USTs at the Site. In a letter dated 23 December 2010, ACEH stated that it would consider case closure of the fuel leak site after the SMP had been implemented and completed, and verification sampling at the base of the excavation at the locations of the former USTs confirmed that petroleum hydrocarbons are below the cleanup levels stated in the SMP. ACEH also requested that groundwater samples be collected from the dewatering system prior to treatment to asses any groundwater contamination. These samples were collected by the dewatering contractor in compliance with NPDES permit requirements.



The general contractor for the Site is DPR Construction, Inc. of Redwood City, California (DPR). The excavation subcontractor for the soil removal and disposal is PSEC. The construction activities involving soil handling began with targeted excavation activities in August 2010 and were completed in May 2011.

Based on information provided to us by DPR and PSEC, approximately 32,020 tons of soil have been excavated and removed from the Site. A total of approximately 17,700 tons of Class II non-hazardous waste was disposed of at Altamont Landfill in Livermore, California, approximately 1,320 tons of Class III non-hazardous waste went to the Dumbarton Quarry (DQA) Landfill in Fremont, California, and approximately 13,000 tons of Class III non-hazardous waste went to Curtner Quarry Landfill in Milpitas, California.

Soil Analytical Results

As stated in the approved SMP, confirmation testing of sidewalls and floor of the final excavation was required to establish the post-excavation soil conditions that would remain beneath the concrete mat foundation and behind the concrete foundation walls. Soil samples were collected with a drive sampler containing $2'' \times 6''$ stainless steel tubes, capped with TeflonTM sheeting and plastic caps, labeled, and placed in an ice-chilled cooler. Samples were submitted to McCampbell Analytical, Inc., a California-certified analytical laboratory in Pittsburg, California under chain-of-custody protocol and documentation.

In accordance with the SMP, the sidewall and bottom samples were analyzed for total petroleum hydrocarbons as gasoline (TPHg) by EPA Method 8015M, TPH as diesel (TPH-d) and as motor oil (TPH-mo) by EPA Method 8015M with silica gel cleanup by EPA Method 3630C, semi-volatile organic compounds (SVOCs) by EPA Method 8270, and total lead. The bottom samples collected beneath the former USTs were also analyzed for benzene, toluene, ethyl benzene, xylenes (BTEX), and methyl tertbutyl ether (MTBE) by EPA Method 8015Bm.

A total of thirty sidewall samples were collected at the Site with fifteen sidewall samples (P5-1 through P5-15) collected at 5 feet below the ground surface (bgs) and fifteen sidewall samples (P10-1 through P10-15) collected at 10 feet bgs between the period of 5 April 2011 through 18 April 2011. A total of sixteen bottom samples were collected at approximately 15 feet bgs (B-1-15 through B-16-15) and one sample at the base of the elevator pit at approximately 21 feet bgs (EP-1-21). Based on analytical results of two of the soil samples (B-1-15 and B-15-15), over-excavation was performed in those areas and additional samples were collected at approximate depths of 17 feet and 17.5 feet bgs (B-1-17 and B-15-17.5) between the period of 15 April 2011 through 28 April 2011.

The analytical results are presented in Tables 1 through 3 and the certified analytical laboratory reports and chain-of-custody records are presented in Appendix A. The soil sample locations are shown on Figure 2.

Residential Screening Levels (SLs) (residential California Human Health Screening Level [CHSSL] for lead) are used as the cleanup goals for soil in the Site to the maximum depth of the redevelopment excavation within the area of the planned garage. The cleanup levels established for soil at the Site are: TPHg and TPHd at 83 milligrams per kilograms (mg/kg), TPHmo at 370 mg/kg, lead at 80 mg/kg, and SVOCs cleanup levels vary per compound.



The soil analytical results are presented in Tables 1 and 2. Analytical results of the soil sample collected beneath the former USTs located in the northern area of the Site that was removed in 1989 (B-1-15) detected TPHg at a concentration of 5.2 mg/kg, TPHd at a concentration of 210 mg/kg, THPmo at a concentration of 160 mg/kg, and total lead at a concentration of 6.6 mg/kg. No BTEX or MTBE were detected at or above the method reporting limit. As the TPHd result was greater than the cleanup level of 83 mg/kg, over-excavation was performed in an area approximately 15 feet by 15 feet by 2 feet in depth. An additional bottom sample was collected at a depth of approximately 17 feet bgs. Bottom sample B-1-17 did not detect TPHg, TPHd, TPHmo, BTEX or MTBE at or above method reporting limits. Total lead was detected at a concentration of 5.2 mg/kg.

Analytical results of the soil sample collected beneath the former UST in the center-eastern area of the Site and removed in 2010 (B-11-15) did not detect TPHg, TPHd, TPHmo, BTEX or MTBE at or above method reporting limits. Total lead was detected at a concentration of 7.3 mg/kg.

Groundwater Analytical Results

The Site currently has a National Pollutant Discharge Elimination System (NPDES) permit from the San Francisco Bay Regional Water Quality Control Board (RWQCB) in place and is operating a dewatering system. The dewatering system consists of 20 perimeter dewatering wells which are pumped through a collection system to one 20,000-gallon baffled settling tank to remove suspended solids. Water is then pumped through two 1,000-pound activated carbon adsorption units arranged in series prior to discharge. The settling tank and carbon units are located on Hollis Street adjacent to the Site. Per the NPDES permit, sampling of the influent and effluent is analyzed for petroleum hydrocarbons, VOCs, SVOCs, metals, pH, total dissolved solids, and turbidity, initially on a monthly basis and subsequently reduced to quarterly. Brighton Environmental Consulting of Oakland, California has been retained by PSEC to oversee and perform the compliance sampling and testing per the NPDES permit requirements. A copy of the March 2010 NPDES Self-Monitoring Report is provided in Appendix B.

The groundwater analytical results are presented in Table 3. As of 26 April 2011, a total of four groundwater samples were collected from the Site's dewatering system prior to any treatment. Analytical results of the initial water sample (DW-1-2/8/11) detected TPHg at a concentration of 63 ug/L, TPHd at a concentration of 97 ug/L, toluene at a concentration of 1.4 ug/L and total xylenes at a concentration of 2.0 ug/L. No benzene, ethylbenzene, or MTBE were detected at or above the method report limits. Analytical results of the water sample INF-001 – 2/28/11 collected by Brighton Environmental Consulting detected TPHmo at a concentration of 51 ug/L and toluene at a concentration of 0.25 ug/L. No TPHg, TPHd, benzene, ethylbenzene, total xylenes, or MTBE were detected at or above method reporting limits. TPHd was the only constituent detected at or above the method reporting limits in Brighton Environmental Consulting's sample INF-001- 3/7/11 at a concentration of 24 ug/L. Treadwell & Rollo collected an additional groundwater sample on 26 April 2011 (DW-1-4/26/11) and no petroleum constituents were detected at or above the method reporting limits.

In our opinion, based on the influent results, any contaminants potentially present in association with releases from the historic USTs has been removed and treated by the construction dewatering system. During the mass excavation activities, residual petroleum hydrocarbons in the soil that could have recontaminated groundwater following cessation of dewatering system have been removed.



Summary

The results of soil confirmation sampling satisfy the established remedial goals and as such the Site is a candidate for case closure. We understand that obtaining an official case closure letter with take some time, so we are requesting communication that at this time we have completed the requirements for case closure. The slab will be poured in less than two weeks. In light of this and in recognition of your administrative process requirements, it would be best for us to speak by telephone to affirm that we have provided all you require for your closure review. I will contact you tomorrow to set up a time for a call next week.

Thanks for all your help with this important redevelopment project.

If you have any questions or require additional information, please contact Peter J. Cusack at 415-955-5244 or by email at: pjcusack@treadwellrollo.com.

Attachments: Tables

Figures

Appendix A - Certified Analytical Laboratory Reports and Chain-of-Custody Records

Appendix B – March 2010 NPDES Self-Monitoring Report

730482302.02 PJC



TABLES

Table 1 Confirmation Soil Samples Analytical Results for Petroleum Hydrocarbons EmeryStation Greenway 5812 Hollis Street

Emeryvile, California Project: 730482302

Sample	Depth	Date	TPHg	TPHd	TPHmo	MTBE	Benzene	Toluene	Ethylbenzene	Xylenes	SVOC
ID Sidewall Sa	feet	Sample					mg,	кg			
P5-1	5	04/05/11	< 1.0	< 1.0	< 5.0						ND
P5-2	5	04/05/11	< 1.0	< 1.0	< 5.0						ND
P5-3	5	04/05/11	< 1.0	< 1.0	< 5.0						ND
P5-4	5	04/05/11	< 1.0	< 1.0	< 5.0						ND
P5-5	5	04/05/11	< 1.0	< 1.0	< 5.0						ND
P5-6	5	04/05/11	< 1.0	< 1.0	< 5.0						ND
P5-7	5	04/05/11	< 1.0	< 1.0	< 5.0						ND
P5-8	5	04/05/11	< 1.0	1.6	6.9						ND
P5-9	5	04/05/11	< 1.0	6.5	14						ND
P5-10	5	04/05/11	< 1.0	8.9	11						1
P5-11	5	04/05/11	580	13,000	5,500						ND
P5-12	5	04/05/11	< 1.0	4.0	< 5.0						ND
P5-13	5	04/05/11	< 1.0	2.4	< 5.0						ND
P5-14	5	04/05/11	< 1.0	< 1.0	< 5.0						ND
P5-15	5	04/05/11	< 1.0	< 1.0	< 5.0						ND
P10-1	10	04/05/11	< 1.0	< 1.0	< 5.0						ND
P10-2	10	04/05/11	< 1.0	< 1.0	< 5.0						ND
P10-3	10	04/05/11	< 1.0	< 1.0	< 5.0						ND
P10-4	10	04/05/11	< 1.0	< 1.0	< 5.0						ND
P10-5	10	04/05/11	< 1.0	< 1.0	< 5.0						ND
P10-6	10	04/05/11	< 1.0	< 1.0	< 5.0						ND
P10-7	10	04/05/11	< 1.0	< 1.0	< 5.0						ND
P10-8	10	04/18/11	< 1.0	1.8	< 5.0						ND
P10-9	10	04/18/11	6.2	100	52						ND
P10-10	10	04/18/11	< 1.0	2.5	< 5.0						ND
P10-11	10	04/15/11	< 1.0	< 1.0	< 5.0						ND
P10-12	10	04/15/11	14	28	13						ND
P10-13	10	04/08/11	< 1.0	< 1.0	< 5.0						ND
P10-14	10	04/08/11	< 1.0	< 1.0	< 5.0						ND
P10-15	10	04/05/11	3.8	370	160						ND
Bottom San											
B-1-15	15	04/15/11	5.2	210	160	< 5.0	< 0.5	< 0.5	< 0.5	< 0.5	ND
B-1-17	17	04/26/11	< 1.0	< 1.0	< 5.0	< 0.05	< 0.005	< 0.005	< 0.005	< 0.005	ND
B-2-15	15	04/15/11	< 1.0	< 1.0	< 5.0						ND
B-3-15	15	04/15/11	< 1.0	< 1.0	< 5.0						ND
B-4-15	15	04/15/11	< 1.0	< 1.0	< 5.0						ND
B-5-15	15	04/15/11	< 1.0	< 1.0	< 5.0						ND
B-6-15	15	04/15/11	< 1.0	< 1.0	< 5.0						ND
B-7-15	15	04/15/11	4.1	45	120						ND
B-8-15	15	04/15/11	< 1.0	< 1.0	< 5.0						ND
B-9-15	15	04/15/11	< 1.0	< 1.0	< 5.0						ND
B-10-15	15	04/18/11	< 1.0	< 1.0	< 5.0						ND
B-11-15	15	04/18/11	< 1.0	< 1.0	< 5.0	< 0.05	< 0.005	< 0.005	< 0.005	< 0.005	ND
B-12-15	15	04/18/11	< 1.0	1.8	< 5.0						ND
B-13-15	15	04/18/11	< 1.0	10	< 5.0						ND
B-14-15	15	04/18/11	3.5	52 630	44						ND
B-15-15	15	04/18/11	21	630	240						ND
B-15-17.5	17.5	04/28/11	79	270	61						ND
B-15-19	19	05/03/11	27	100	31						ND
B-16-15	15 21	04/18/11	< 1.0	< 1.0 < 1.0	< 5.0						ND
EP-1 Cleanup Go (mg/kg) (R Shallow Soi Drinking Wa	als for the esidentia I, Ground	l ESL, water is	< 1.0	83	< 5.0 370	NA	NA	NA	NA	NA	ND Various

Notes:

mg/kg - milligrams per kilograms

TPHg - Total Petroleum Hydrocarbons as Gasoline, EPA Method 8015M TPHd - Total Petroleum Hydrocarbons as Diesel Range, EPA Method 8015M

TPHmo - Total Petroleum Hydrocarbons as Motor Oil, EPA Method 8015M

MTBE - Methyl Tert-Butyl Ether, EPA Method 8260B

Benzene, Toluene, Ethylbenzene, and Xylenes, EPA Method 8260B

 $\ensuremath{\mathsf{SVOC}}$ - semi-volatile organic compounds, EPA Method 8270

Concentration listed in **bold** indicate an exceedance of applicable cleanup levels.

NA - Not Applicable

¹ - **0.42** mg/kg Benzo(a)anthracene, **0.49** mg/kg Benzo(k)fluoranthene, **0.79**mg/kg Beno(a)pyrene, 0.51 mg/kg Chrysene, 1.1 mg/kg Flouranthene, **0.68** mg/kg Indeno(1,2,3-cd)pyrene, **0.60** mg/kg Benzo(b)fluoranthene, 1.2 mg/kg Benzo(g,h,i)perylene, 0.48 mg/kg Phenanthrene, and 1.9 mg/kg Pyrene are detected in P5-10.

Table 2 **Confirmation Soil Samples Analytical Results for Total Lead EmeryStation Greenway 5812 Hollis Street Emeryvile, California** Project: 730482302

Sample ID	Depth interval feet	Date Sampled	Lead
			(mg/kg)
Sidewall Sam	ples		
P5-1	5	04/05/11	7.8
P5-2	5	4/5/2011	8.6
P5-3	5	4/5/2011	< 5.0
P5-4	5	4/5/2011	7.8
P5-5	5	4/5/2011	8.9
P5-6	5	4/5/2011	10
P5-7	5	4/5/2011	5.1
P5-8	5	4/5/2011	14
P5-9	5	4/5/2011	8.6
P5-10	5	4/5/2011	15
P5-11	5	4/5/2011	7.5
P5-12	5	04/05/11	8.3
P5-13	5	04/05/11	6.4
P5-14	5	04/05/11	6.2
P5-15	5	04/05/11	8.4
P10-1	10	04/05/11	9.5
P10-2	10	04/05/11	< 5.0
P10-3	10	04/05/11	7.1
P10-4	10	04/05/11	< 5.0
P10-5	10	04/05/11	< 5.0
P10-6	10	04/05/11	8
P10-7	10	04/05/11	6
P10-8	10	04/18/11	11
P10-9	10	04/18/11	9.9
P10-10	10	04/18/11	12
P10-11	10	04/15/11	13
P10-12	10	04/15/11	5.1
P10-13	10	04/08/11	8.7
P10-14	10	04/08/11	9.3
P10-15	10	04/05/11	6.3
Bottom Samp		2 4 224 ==	
B-1-15	15	04/15/11	6.6
B-1-17	17	04/26/11	5.2
B-2-15	15	04/15/11	5.6
B-3-15	15	04/15/11	8.4
B-3-15 B-4-15	15	04/15/11	12
B-5-15	15	04/15/11	7.8
B-6-15	15	04/15/11	5.7
B-7-15	15	04/15/11	8.6
B-7-15 B-8-15	15	04/15/11	11
B-9-15	15	04/15/11	7.6
B-10-15	15	04/18/11	< 5.0
B-11-15	15	04/18/11	7.3
B-12-15	15	04/18/11	8.2
B-13-15	15	04/18/11	13
B-14-15	15	04/18/11	7.6
B-15-15	15	04/18/11	7.2
B-15-17.5	17.5	04/28/11	9.1
B-16-15	15	04/18/11	7.0
EP-1 Cleanup Goal (Residential (-		8.6 80

Notes:

mg/kg - milligrams per kilograms < 5.0 - Analyte was not detected above the laboratory reporting limit (5.0 mg/kg).

Table 3 Confirmation Influent Groundwater Samples Analytical Results for Petroleum Hydrocarbons EmeryStation Greenway 5812 Hollis Street

Emeryvile, California Project: 730482302

Sample ID	Date Sampled	TPHg	TPHmo	TPHd	Benzene	ethylbenzene	Toluene	Total Xylenes	МТВЕ
	Jampiou					μg/L			
DW-1	02/08/11	63		97	< 0.5	< 0.5	1.4	2.0	< 0.5
INF-001	02/28/11	< 51	51 J B	< 52	< 0.5	< 0.5	0.25 J	< 1.0	< 0.5
INF-001	03/07/11	< 51	< 100	24 J B	< 0.5	< 0.5	< 0.5	< 1.0	< 0.5
DW-1	04/26/11	< 50	< 250	< 50	< 0.5	< 0.5	< 0.5	< 0.5	< 0.5

Notes:

All results are reported in microgram per liter (µg/L)

TPHg - Total Petroleum Hydrocarbons as Gasoline, EPA Method 8015M

TPHmo - Total Petroleum Hydrocarbons as Moto OII Range (C10-C23), EPA Method 8015M

TPHd - Total Petroleum Hydrocarbons as Diesel Range (C10-C23), EPA Method 8015M

MTBE - Metyl Tertiary Butyl

- -- Not Analyzed
- J J-flag by laboratroy, indicating estimated concentration is below reporting limit but above method detection limit
- B compound was found in the blank and sample
- < 0.5 Analyte was not detected above the laboratory reporting limit (0.5 μ g/L)
- ND Not detected at or above the laboratory reporting limit

Table 4 Confirmation Soil Samples Analytical Results for Total Metals EmeryStation Greenway 5812 Hollis Street Emeryvile, California

Project: 730482302

Sample ID	Date Sampled	Antimony	Arsenic	Barium	Beryllium	Cadmium	Chromium	Cobalt	Copper	Lead	Mercury	Molybdenum	Nickel	Selenium	Silver	Thallium	Vanadium	Zinc
10	Sampled									mg/kg	}							
DW-1	02/08/11	< 0.5	< 0.5	< 5.0	< 0.5	< 0.25	< 0.5	< 0.5	< 0.5	< 0.5	< 0.025	< 0.5	< 0.5	< 0.5	< 0.19	< 0.5	< 0.5	< 5.0

Notes

All results are reported in milligrams per kilogram (mg/kg)

-- Not analyzed

< 0.5 - Analyte was not detected above the laboratory reporting limit (0.5 mg/kg)

TTLC - Total Threshold Limited Concentration

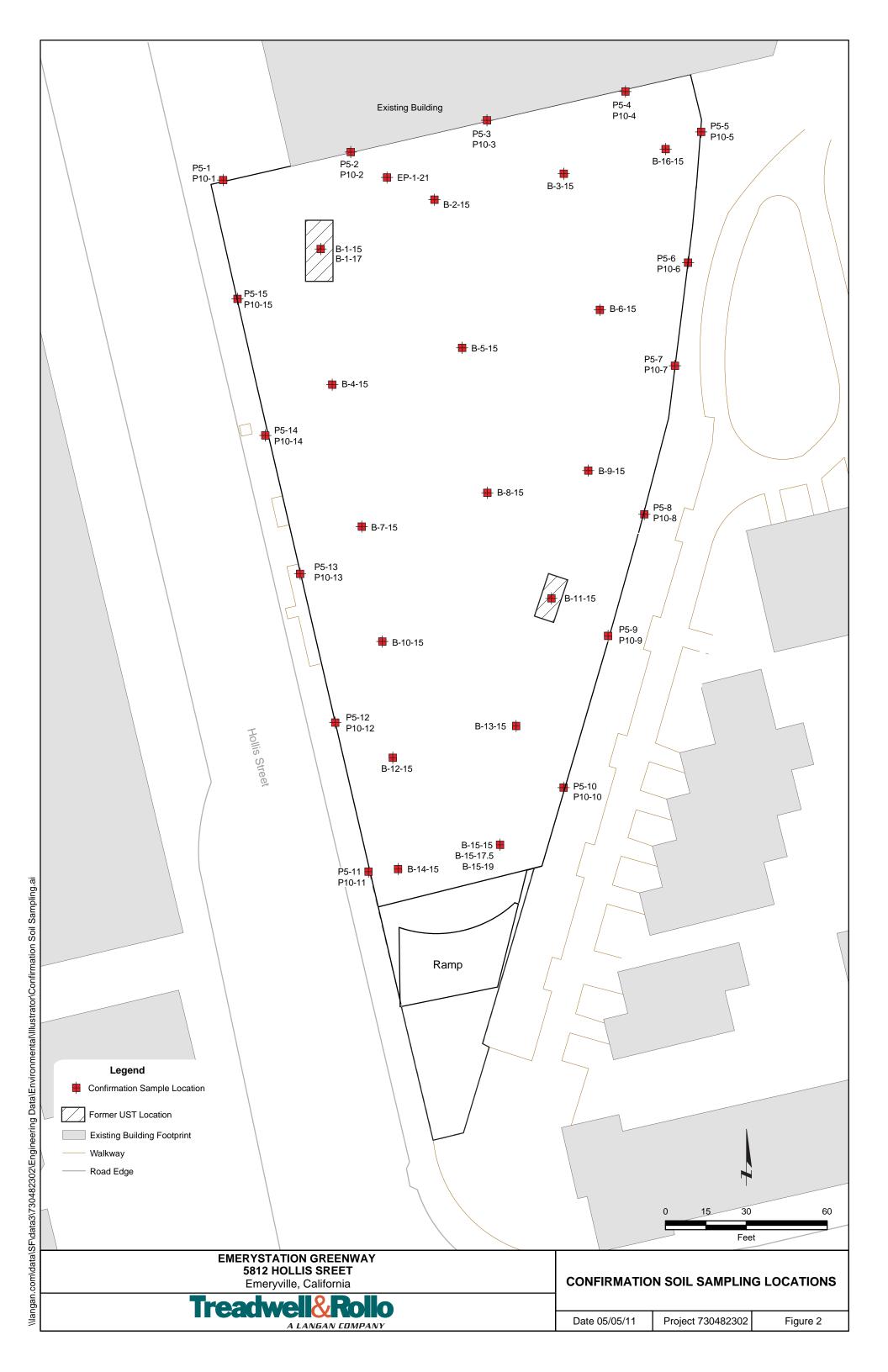
STLC - Soluble Threshold Limited Concentration

TCLP - Toxicity Characteristic Leaching Procedure



FIGURES







APPENDIX A

Certified Analytical Laboratory Reports and Chain-of-Custody Records

McCampbell Analytical, Inc.
"When Quality Counts"

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300		Date Received: 04/07/11
oss frongomery bu, butter 1900	Client Contact: Peter Cusack	Date Reported: 04/13/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/26/11

WorkOrder: 1104217

April 26, 2011

Enclosed within are:

- 1) The results of the 8 analyzed samples from your project: #730482302; 5812 Hollis St,
- 2) A QC report for the above samples,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius
Laboratory Manager
McCompbell Application Inc.

McCampbell Analytical, Inc.

1104217

Treatwe Environmental and Geote				ontg	omer	Str	eet, S	Suite	1300), Sa	n Fra	ncis	co,	CO CA 941 510.87	11 Ph:	: 41						15.9	955.	9041				Pag	e _(of <u></u>
Site Name: Job Number: Project Manager\Cor	() . I	Hallis Return		lec		mmg		ld., S						CA 95		h: 9	16.5	65.	741	2/Fa		916	5.56	5.7412			1	urnar Tim	e ,	
Samplers: Recorder (Signature	Required):	W.Jano	and				No.	Con	tain	ers	10											clean-up					110	M	al	_
, , ,	, , ,	-01		M	atrix	-			ervat		d	7	_									clear								
Field Sample Identification No.	Date	Time	Lab Sample No.	Soil	Water	E CH	H ₂ SO ₄	HNO3	90	Omer	PIK	SVA	Lead									Silica gel	Hold				Rema	arks		
P10-1	4-5-11	5935		X		Т			X.		X	X	X			T	T		T	T		X								
P10-Z	4-5-11	0945		X		Т			X		X	X	X				T	T	T	T	1	X								
P10-3	4-5-11	0950		X	2	I			Ø-	,	X	C	X			I	I	I			1	X								
P10-4	4-5-11	1020		X		1			1	-	X	X	X			1	1	1	1	1	-	X								
P10-5	4-5-11	1025		X	4	+		Ц	X.	+	X	X	X			+	+	+	4	+	4	S								
P10-6	4-3-11	1035		X	/	+	+	\vdash	X)	+	0	X	X		-	+	+	+	+	+	+	X								
P10-7	4-5-11	1045		×	+	+	+	H	A	+	X	×	X	-	-	+	+	+	+	+	-1	^	-							
410-15	4-6-11	1140		Н	+	+	+	H	+	+	+		-		+	+	1	4	1	H	+	\dashv	-							
				Н	+	+	+	Н	+	+	Н	\dashv		-	IC		100	ONE	5171	ON	+	+	۸	PPRO	PRIAT	TE				
				H	+	+	+	H	+	+	+	\dashv	+	+	H	EAL) BP	ACI	EA	BSE	N			CON	TAIN	ERS	N LAE			
				Н	+	$^{+}$	+	H	+	+	Н				D	ECH	140	RN		_	DAS	AB	& G	META			N LAD			
				H	+	$^{+}$	+	\vdash	+	+	\vdash				P	RES	ER	VAT	ICA	+	+	+			_					
				Н	+	$^{+}$	$^{+}$	\forall	+	+		1				†	+	t	,	+	†	\forall			1					
Relinquished by: (Signati	MAA	97	Date 4-77-	11		Ti	me	4	20	5	Red	eive	ed by	: (Sign	Hure)		1	1			-	Date	4	/	7/	1	Time	19.	2	0
Relinquished by: (Silgnate	ure)	1	Date 4/7/1	/		Ti	me/	7	45	5	Red	eive	d by	: (Signa	ature)			1				Date	/	7			Time			
Relinquished by: (Signatu	ure)		Date /			Ti	me				Red	erve	dipy	Lab: (Signatu	ure)	/	2	1	/	[Date	2	47	//	/	Time	71	5	•
Sent to Laboratory (Laboratory Commen		McC	ampbell							_	Me			Ship d Carrie		Pr				couri		Nar		Fed Ex	x		Airbo	rne		UPS
		White Copy	- Original		Yello	w C	ору -	- Lat	borat	tory	_			Pin	k Cop	oy -	Fie	ld				С	00	Nun	nber	0	05	23	7	

1534 Willow Pass Rd Pittsburg, CA 94565-1701 (925) 252-9262

CHAIN-OF-CUSTODY RECORD

San Francisco, CA 94111

Page 1 of 1

Date Printed: 04/08/2011

(925) 252-9262				WorkOr	der: 110421	7 ClientC	ode: TWRF		
	WaterTrax	WriteOn	☐ EDF	Excel	Fax	✓ Email	HardCopy	ThirdParty	☐ J-flag
Report to:				Bil	II to:		Req	uested TAT:	5 days
Peter Cusack	Email: p	pjcusack@treadv	wellrollo.com		Accounts Pa	ayable			
Treadwell & Rollo	cc:				Treadwell &	Rollo	_		
555 Montgomery St., Suite 1300	PO:				555 Montgor	mery St., Suite 130	0 Dat	e Received:	04/07/2011

ProjectNo: #730482302; 5812 Hollis St

(415) 955-5244 FAX (415) 955-9041

San Francisco, CA 94111

					Requested Tests (See legend below)													
Lab ID	Client ID	Matrix	Collection Date	Hold	1	2	3	4	5	6	7	8	9	10	11	12		
1104217-001	P10-1	Soil	4/5/2011 9:35		Α	Α	Α											
1104217-002	P10-2	Soil	4/5/2011 9:45		Α	Α	Α											
1104217-003	P10-3	Soil	4/5/2011 9:50		Α	Α	Α											
1104217-004	P10-4	Soil	4/5/2011 10:20		Α	Α	Α											
1104217-005	P10-5	Soil	4/5/2011 10:25		Α	Α	Α											
1104217-006	P10-6	Soil	4/5/2011 10:35		Α	Α	Α											
1104217-007	P10-7	Soil	4/5/2011 10:45		Α	Α	Α											
1104217-008	P10-15	Soil	4/5/2011 11:40		Α	Α	Α											

Test Legend:				
1 8270D_S	2 PB_S	3 TPH(D)WSG_S	4	5
6	7	8	9	10
11	12			
The following SampIDs: 001A, 002/	A, 003A, 004A, 005A, 006A, 007A, 008A cor	ntain testgroup.		Prepared by: Ana Venegas

Comments: SEND HARD COPY

NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days).

Hazardous samples will be returned to client or disposed of at client expense.

Sample Receipt Checklist

Client Name:	Treadwell & R	ollo			Date a	and Time Received:	4/7/2011 7	7:25:54 PM
Project Name:	#730482302; 5	812 Hollis St			Check	list completed and r	eviewed by:	Ana Venegas
WorkOrder N°:	1104217	Matrix Soil			Carrie	r: Rob Pringle (M	IAI Courier)	
		<u>Chair</u>	ı of Cu	stody (C	COC) Informa	ition		
Chain of custody	y present?		Yes	V	No 🗆			
Chain of custody	y signed when relir	nquished and received?	Yes	V	No 🗆			
Chain of custody	y agrees with sam	ole labels?	Yes	✓	No 🗆			
Sample IDs noted	d by Client on COC	?	Yes	V	No 🗆			
Date and Time o	f collection noted b	y Client on COC?	Yes	~	No 🗆			
Sampler's name	noted on COC?		Yes	V	No 🗆			
		<u>s</u>	ample	Receipt	t Information			
Custody seals in	ntact on shipping co	ontainer/cooler?	Yes		No 🗆		NA 🗹	
Shipping contain	ner/cooler in good o	ondition?	Yes	V	No 🗆			
Samples in prop	er containers/bottl	es?	Yes	V	No 🗆			
Sample containe	ers intact?		Yes	✓	No 🗆			
Sufficient sample	e volume for indica	ted test?	Yes	✓	No 🗌			
		Sample Prese	rvatio	n and Ho	old Time (HT)) Information		
All samples rece	eived within holding	time?	Yes	✓	No 🗌			
Container/Temp	Blank temperature		Coole	er Temp:	10.4°C		NA \square	
Water - VOA via	als have zero head	space / no bubbles?	Yes		No 🗆	No VOA vials subm	itted 🗹	
Sample labels cl	hecked for correct	preservation?	Yes	~	No 🗌			
Metal - pH accep	otable upon receipt	(pH<2)?	Yes		No 🗆		NA 🔽	
Samples Receive	ed on Ice?		Yes	V	No 🗆			
		(Ice Typ	e: WE	TICE)			
* NOTE: If the "I	No" box is checke	d, see comments below.						
=====	=====	======			====	=====	====	
Client contacted:	:	Date contac	ted:			Contacted	by:	
Comments:								

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104217

Lab ID		1104217-001A					
Client ID		P10-1					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surro	gate Re	coveries (%)			
%SS1:	102	2		%SS2:	102	2	
0.000				0.004			

Surrogate Recoveries (%) %SS1: 102 %SS2: 102 %SS3: 99 %SS4: 83 %SS5: 79 %SS6: 86

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	3	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Work Order: 1104217 Extraction Method: SW3550B

Lab ID		1104217-002A					
Client ID		P10-2					
Matrix		Soil					
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33	, , , , , , , , , , , , , , , , , , , ,	·		
		Surro		coveries (%)			
%SS1:	103		g	%SS2:	104	1	
%SS3:	100			%SS4:	86		

Surrogate Recoveries (%)					
%SS1:	103	%SS2:	104		
%SS3:	100	%SS4:	86		
%SS5:	84	%SS6:	89		
	·	·			

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104217

Lab ID		1104217-003A					
Client ID		P10-3					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surro	gate Re	coveries (%)			•
%SS1:	100)		%SS2:	101	1	

Surrogate Recoveries (%)					
%SS1:	100	% SS2:	101		
%SS3:	97	%SS4:	84		
%SS5:	75	%SS6:	85		

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	3	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104217

Lab ID		1104217-004A					
Client ID		P10-4					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surro	gate Re	coveries (%)			
%SS1:	100)		%SS2:	100)	
** 000	100						

Surrogate Recoveries (%)							
%SS1:	100	%SS2:	100				
%SS3:	98	%SS4:	83				
%SS5:	76	%SS6:	86				

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in $\mu g/L$, soil/sludge/solid samples in mg/kg, wipe samples in $\mu g/wipe$, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104217

Lab ID		1104217-005A					
Client ID				P10-5			
Matrix		Soil					
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surre	gate Re	coveries (%)			
%SS1:	103			%SS2:	102	2	

Surrogate Recoveries (%)					
%SS1:	103	% SS2:	102		
%SS3:	100	%SS4:	86		
%SS5:	75	%SS6:	87		
		· ·			

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	3	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104217

Lab ID		1104217-006A					
Client ID		P10-6					
Matrix				Soil			
Compound	Concentration *	on * DF Reporting Limit Compound		Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
	-	Surre	gate Re	coveries (%)		·	·
0/ 001.	102		-	0/ 992	100		

Surrogate Recoveries (%)					
%SS1:	102	%SS2:	100		
%SS3:	99	%SS4:	85		
%SS5:	74	%SS6:	87		

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in µg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

McCampbell Analytical, Inc. "When Quality Counts"

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

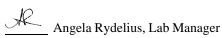
Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104217

Lab ID	1104217-007A						
Client ID				P10-7			
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
	-	Surre	ogate Re	coveries (%)	-	· · ·	
%SS1:	105	;		%SS2:	103	3	
% SS3·	103			% SSA·	86		

%SS1: 105 %SS2: 103 %SS3: 102 %SS4: 86 %SS5: 75 %SS6: 88

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	3	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104217

Lab ID	1104217-008A						
Client ID		P10-15					
Matrix		Soil					
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33	·			
		Surro	gate Re	coveries (%)			
%SS1:	103			%SS2:	96		
0/ 552	102			0/ 554.	02		

Surrogate Recoveries (%) %SS1: 103 %SS2: 96 %SS3: 104 %SS4: 83 %SS5: 68 %SS6: 84

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
<i>5</i> ,	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 04/08/11-04/12/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline*

Extraction method	SW5030B	Analytical methods SW8015Bm		Work Order:		1104217
Lab ID	Client ID	Matrix	TPH(g)	DF	% SS	Comments
001A	P10-1	S	ND	1	74	
002A	P10-2	S	ND	1	82	
003A	P10-3	S	ND	1	86	
004A	P10-4	S	ND	1	88	
005A	P10-5	S	ND	1	70	
006A	P10-6	S	ND	1	78	
007A	P10-7	S	ND	1	83	
008A	P10-15	S	3.8	1	79	d7
	porting Limit for DF =1;	W	NA		NA	
	means not detected at or cove the reporting limit	S	1.0		mg/Kg	3

* water and vapor samples are reported in µg/L, soil/sludge/solid samples in mg/kg,	wipe samples in µg/wipe, product/oil/non-aqueous liquid
samples and all TCLP & SPLP extracts in mg/L	

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

d7) strongly aged gasoline or diesel range compounds are significant in the TPH(g) chromatogram

Angela Rydelius, Lab Manager

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:



Extraction method: SW3050B

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/08/11

Lead by ICP*

Analytical methods: SW6010B

Lab ID Client ID Matrix Extraction Type Lead DF % SS Comments 1104217-001A S TOTAL 9.5 1 97 P10-1 1104217-002A P10-2 S TOTAL ND 1 96 1104217-003A S TOTAL P10-3 7.1 1 92 1104217-004A P10-4 S TOTAL ND 1 98 1104217-005A P10-5 S TOTAL ND 1104217-006A P10-6 S TOTAL 8.0 92 1104217-007A P10-7 S TOTAL 1 94 6.0 1104217-008A S TOTAL P10-15 1 89 6.3

Reporting Limit for DF =1;	W	TOTAL	NA	μg/L
ND means not detected at or above the reporting limit	S	TOTAL	5.0	mg/Kg

*water samples are reported in μ g/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, filter samples in μ g/filter.

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of $0.45 \mu m$ filtered and acidified sample.

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

Angela Rydelius, Lab Manager

DHS ELAP Certification 1644

Work Order: 1104217



	Client Project ID: #730482302; 5812	Date Sampled:	04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received:	04/07/11
	Client Contact: Peter Cusack	Date Extracted:	04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/08/11-04/12/11

Total Extractable Petroleum Hydrocarbons with Silica Gel Clean-Up*

Extraction method: SW3550B/3630C Analytical methods: SW8015B Work Order: 1104217 TPH-Diesel TPH-Motor Oil % SS Lab ID Client ID DF Matrix Comments (C10-C23) (C18-C36) 1104217-001A P10-1 S ND ND 104 1104217-002A P10-2 S ND ND 92 1104217-003A P10-3 S ND ND 104 1104217-004A P10-4 S ND ND 107 1104217-005A P10-5 S ND ND 107 1 1104217-006A S 108 P10-6 ND ND 1104217-007A S ND 108 P10-7 ND 1 1104217-008A P10-15 S 370 160 1 107 e3,e7

Reporting Limit for DF =1;	W	NA	NA	ug/L
ND means not detected at or above the reporting limit	S	1.0	5.0	mg/Kg

^{*} water samples are reported in μ g/L, wipe samples in μ g/wipe, soil/solid/sludge samples in mg/kg, product/oil/non-aqueous liquid samples in mg/L, and all DISTLC / STLC / SPLP / TCLP extracts are reported in μ g/L.

cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract.

%SS = Percent Recovery of Surrogate Standard. DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

e3) aged diesel is significant

e7) oil range compounds are significant



QC SUMMARY REPORT FOR SW8270C

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57506 WorkOrder 1104217

EPA Method SW8270C	Extra	ction SW	3550B					S	Spiked San	nple ID	: 1104143-0	01A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%))
Analyte	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Acenaphthene	ND	2	61.2	64.9	5.81	64.8	65.1	0.431	30 - 130	30	30 - 130	30
4-Chloro-3-methylphenol	ND	4	59.8	63.7	6.32	72	70.3	2.34	30 - 130	30	30 - 130	30
2-Chlorophenol	ND	4	60.8	57.1	6.31	61.7	62.9	1.97	30 - 130	30	30 - 130	30
1,4-Dichlorobenzene	ND	2	66.4	70.3	5.76	68.3	67.9	0.543	30 - 130	30	30 - 130	30
2,4-Dinitrotoluene	ND	2	61.2	68.1	10.7	74.3	72.5	2.44	30 - 130	30	30 - 130	30
4-Nitrophenol	ND	4	89.2	89.6	0.492	72.1	67.9	5.98	30 - 130	30	30 - 130	30
N-Nitrosodi-n-propylamine	ND	2	59.8	53.4	11.4	55.6	54.8	1.47	30 - 130	30	30 - 130	30
Pentachlorophenol	ND	4	54	53.9	0.297	46.6	45.6	2.30	30 - 130	30	30 - 130	30
Phenol	ND	4	61.4	59.8	2.65	63.3	62	2.02	30 - 130	30	30 - 130	30
Pyrene	ND	2	49.7	50.2	1.00	62.5	54.6	13.6	30 - 130	30	30 - 130	30
1,2,4-Trichlorobenzene	ND	2	74.7	73.9	1.02	79.3	77.3	2.53	30 - 130	30	30 - 130	30
%SS1:	81	200	76	73	4.02	84	83	1.80	30 - 130	30	30 - 130	30
%SS2:	77	200	72	73	0.730	89	87	1.85	30 - 130	30	30 - 130	30
%SS3:	82	200	83	79	4.30	96	94	2.41	30 - 130	30	30 - 130	30
%SS4:	80	200	80	80	0	88	89	0.587	30 - 130	30	30 - 130	30
%SS5:	67	200	71	66	6.85	98	95	3.22	30 - 130	30	30 - 130	30
%SS6:	74	200	74	78	5.72	104	94	9.34	30 - 130	30	30 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57506 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104217-001A	04/05/11 9:35 AM	04/07/11	04/09/11 9:09 AM	1104217-002A	04/05/11 9:45 AM	04/07/11	04/09/11 10:33 AM
1104217-003A	04/05/11 9:50 AM	04/07/11	04/09/11 11:57 AM	1104217-004A	04/05/11 10:20 AM	04/07/11	04/09/11 1:25 PM
1104217-005A	04/05/11 10:25 AM	04/07/11	04/09/11 2:52 PM	1104217-006A	04/05/11 10:35 AM	04/07/11	04/09/11 4:22 PM
1104217-007A	04/05/11 10:45 AM	04/07/11	04/09/11 5:46 PM	1104217-008A	04/05/11 11:40 AM	04/07/11	04/09/11 7:06 PM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and / or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix, sample diluted due to high matrix or analyte content, or MS/MSD samples diluted due to high organic content.

#) surrogate diluted out of range; & = low or no recovery of surrogate or target analytes due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.

QA/QC Officer

QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57512 WorkOrder 1104217

EPA Method SW8015B	S	piked San	nple ID:	: 1104154-0	002A							
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
, and y to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	2.5	40	119	118	1.02	115	113	2.17	70 - 130	30	70 - 130	30
%SS:	106	25	108	107	0.632	95	93	2.07	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57512 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104217-001A	04/05/11 9:35 AM	04/07/11	04/11/11 8:00 PM	1104217-002A	04/05/11 9:45 AM	04/07/11	04/12/11 6:01 AM
1104217-003A	04/05/11 9:50 AM	04/07/11	04/11/11 6:51 PM	1104217-004A	04/05/11 10:20 AM	04/07/11	04/09/11 2:00 PM
1104217-005A	04/05/11 10:25 AM	04/07/11	04/08/11 4:08 PM	1104217-006A	04/05/11 10:35 AM	04/07/11	04/08/11 7:43 AM
1104217-007A	04/05/11 10:45 AM	04/07/11	04/08/11 7:45 PM	1104217-008A	04/05/11 11:40 AM	04/07/11	04/08/11 5:28 AM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

DHS ELAP Certification 1644

QA/QC Officer

QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57513 WorkOrder 1104217

EPA Method SW8015Bm	Extrac	ction SW	5030B					5	Spiked Sar	nple ID	: 1104152-0	003A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acc	eptance	Criteria (%)	1
7 tildiyto	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex)	ND	0.60	124	125	0.352	128	124	2.79	70 - 130	20	70 - 130	20
MTBE	ND	0.10	101	109	7.37	108	112	3.44	70 - 130	20	70 - 130	20
Benzene	ND	0.10	89.5	90.4	1.03	91.1	93.6	2.67	70 - 130	20	70 - 130	20
Toluene	ND	0.10	88.1	88.8	0.737	88.7	91.8	3.39	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	89.5	90.6	1.24	90.8	94.3	3.81	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	89.5	90.9	1.51	90.6	93.9	3.57	70 - 130	20	70 - 130	20
%SS:	83	0.10	78	82	4.96	76	87	12.8	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57513 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104217-001A	04/05/11 9:35 AM	04/07/11	04/12/11 1:47 AM	1104217-002A	04/05/11 9:45 AM	04/07/11	04/08/11 6:34 PM
1104217-003A	04/05/11 9:50 AM	04/07/11	04/08/11 7:04 PM	1104217-004A	04/05/11 10:20 AM	04/07/11	04/08/11 7:34 PM
1104217-005A	04/05/11 10:25 AM	04/07/11	04/12/11 2:46 AM	1104217-006A	04/05/11 10:35 AM	04/07/11	04/12/11 2:16 AM
1104217-007A	04/05/11 10:45 AM	04/07/11	04/08/11 9:04 PM	1104217-008A	04/05/11 11:40 AM	04/07/11	04/09/11 6:54 AM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

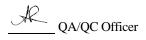
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104217

EPA Method SW6010B		Extraction SW3050B				BatchID: 57460			Spike	Spiked Sample ID: 1104093-003A			
Analyte	Sample	Spiked	MS	MSD	MS-MSD	Spiked	LCS	LCSD	LCS-LCSD	Acc	eptanc	e Criteria (%	5)
Analyte	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Lead	15	50	96.9	104	5.61	10	101	95.6	5.17	75 - 125	25	75 - 125	25
%SS:	98	500	98	95	2.64	500	103	100	3.84	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57460 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracte	d Date Analyzed
1104217-001A	04/05/11 9:35 AM	04/07/11	04/08/11 2:49 PM	1104217-002A	04/05/11 9:45 AM	04/07/11	04/08/11 2:51 PM
1104217-003A	04/05/11 9:50 AM	04/07/11	04/08/11 2:53 PM	1104217-004A	04/05/11 10:20 AM	04/07/11	04/08/11 2:56 PM
1104217-005A	04/05/11 10:25 AM	04/07/11	04/08/11 3:02 PM	1104217-006A	04/05/11 10:35 AM	04/07/11	04/08/11 3:05 PM
1104217-007A	04/05/11 10:45 AM	04/07/11	04/08/11 3:07 PM	1104217-008A	04/05/11 11:40 AM	04/07/11	04/08/11 3:09 PM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer

McCampbell Analytical,	Inc.
"XXII O1' C	

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300		Date Received: 04/07/11
oss frongomery bu, butter 1900	Client Contact: Peter Cusack	Date Reported: 04/14/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/26/11

WorkOrder: 1104220

April 26, 2011

-	-
llaar	Peter:
17541	FEIGI.

Enclosed within are:

- 1) The results of the 15 analyzed samples from your project: #730482302; 5812 Hollis St,
- 2) A QC report for the above samples,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius
Laboratory Manager
McGamphall Application Local

McCampbell Analytical, Inc.



Treadwe			CHA S 555 M													55.9040	/Fax:	415.9	55.90	41		Page		2
Site Name: Job Number:	5812	B 4.	S St Alds	4th Stre	et, T	hird	Floor	, Oak	land (CA 94	612	Ph: 510 nto, CA	0.874 A 9582	.4500/ 25 Ph	/Fax: 5	510.874. 565.741 uestec	.4507 2/Fax					Turnar	ound	
Project Manager\Cor Samplers: Recorder (Signature	Rob	Mila		Mat	rix		lo. Co			10 -d	136							clean-up				Vol	Ma	(
Field Sample Identification No.	Date	Time	Lab Sample No.	Soil	Other	HCL	H ₂ SO ₄	lce	Other	TPH	SVOC	lend			Ш	Ш			Hold		Re	marks		
P5-1	4-5-11	1351		X				X		X	X	X		+			+	X	+					
95-4 95-4	4-4-11	1334		X				X		X	X	X.		+	H	+	+	X	+					
P5-6	4-5-11	1030		X				X		X	X	X						X X						
P5-9	4-5-11	0925		X				X		X	XXX	X		+		+	+	XXX	+					=
P5-12	4-5-11	0915		X X				XX		X	XX	X				\pm		X	\pm					
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Sent to Laboratory (Name): Laboratory Comments/Notes:												Method of Shipment Lab courier Fed Ex Airborne Hand Carried Private Courier (Co. Name)											UP	S
		White Copy	- Original	Ye	ellow	Co	py - La	abor	atory	_			Pink	Сор	HE	OD CON AD SPAC	CEAB	SENT DIN L	AB_	Number: APPROPR CONTA PRESI	ERVED	IN LAB_	6	

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Site Name: Job Number: Project Manager\Cor Samplers: Recorder (Signature	Rob	Hollis 82307 Helan	4 Aide	K)ap -	No.		aine	ers	20.00	्र रा	- No.	Analy:					clean-up	.56:	5.7412	Turnaround Time	
Field Sample Identification No.	Date	Time	Lab Sample No.	Soil	Water	크	$\overline{}$	_	lce	$\overline{}$	161	SVO	Lead		Ш				Silica gel	Hold		Remarks	
P5-15	4-5-11	0940										<u>X</u>	X						×				
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McCampbell Analytical, Inc.

1534 Willow Pass Rd

CHAIN-OF-CUSTODY RECORD

Page 1 of 2

Pittsburg (925) 25	g, CA 94565-1701 52-9262					Work	Order	: 1104	220		Client	Code: T	WRF				
		WaterTrax	WriteOn	☐ EDF		Excel		Fax		✓ Emai	I	Hard	Сору	Thi	rdParty	☐ J-	flag
-	Rollo mery St., Suite 1300 co, CA 94111	cc: PO:		adwellrollo.com 5812 Hollis St			Tre 55		l & Ro gomer			00	Dat	uested e Rece e Prin	ived:	5 0 04/07/ 04/08/	
									Re	auestea	l Tests	(See le	gend b	elow)			
Lab ID	Client ID		Matrix	Collection Date	Hold	1	2	3	4	5	6	7	8	9	10	11	12
1104220-001	P5-1		Soil	4/5/2011 9:30		Α	Α	Α									
1104220-002	P5-2		Soil	4/5/2011 13:21		Α	Α	Α									
1104220-003	P5-3		Soil	4/5/2011 13:24		Α	Α	Α									
1104220-004	P5-4		Soil	4/5/2011 13:30		Α	Α	Α									
1104220-005	P5-5		Soil	4/5/2011 13:35		Α	Α	Α									
1104220-006	P5-6		Soil	4/5/2011 10:30		Α	Α	Α									
1104220-007	P5-7		Soil	4/5/2011 10:40		Α	Α	Α									
1104220-008	P5-8		Soil	4/5/2011 10:50		Α	Α	Α									
1104220-009	P5-9		Soil	4/5/2011 9:25		Α	Α	Α									
1104220-010	P5-10		Soil	4/5/2011 9:20		Α	Α	Α									
1104220-011	P5-11		Soil	4/5/2011 9:15		Α	Α	Α									
1104220-012	P5-12		Soil	4/5/2011 9:10		Α	Α	Α									
1104220-013	P5-13		Soil	4/5/2011 9:05		Α	Α	Α									
1104220-014	P5-14		Soil	4/5/2011 13:46		Α	Α	Α									
Test Legend:																	
1 8270	DD_S 2	PB_S		3 TP	H(D)W	SG_S		4	ı					5			
6	7			8				9)					10			
11	12																
The following San testgroup.	npIDs: 001A, 002A, 003A, 00	04A, 005A, 006A	, 007A, 008A,	009A, 010A, 011A,	012A,	013A,	014A, 0	15A cor	ntain				Prep	ared by	: Ana	Venegas	<u>s</u>

Comments: SEND HARD COPY

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

McCampbell Analytical, Inc.

1534 Willow Pass Rd

CHAIN-OF-CUSTODY RECORD

Page 2 of 2

Pittsburg, CA (925) 252-92						Work(Ordei	:: 1104	220	(ClientC	ode: T	WRF				
		WaterTrax	WriteOn	☐ EDF		Excel		Fax		Email		Hard	Сору	Third	lParty	☐ J-f	lag
Report to:							Bill to	:					Req	uested 1	TAT:	5 d	lays
Peter Cusack Treadwell & Roll 555 Montgomery San Francisco, C (415) 955-5244	St., Suite 1300	cc: PO:		adwellrollo.com 5812 Hollis St			T:		I & Roll gomery)		e Recei e Printe		04/07/2 04/08/2	
									Req	uested	Tests (See leg	end b	elow)			
Lab ID	Client ID		Matrix	Collection Date	Hold	1	2	3	4	5	6	7	8	9	10	11	12
Lab ID 1104220-015	Client ID P5-15		Matrix Soil	4/5/2011 9:40	Hold	1	2	3 A	4	5	6	7	8	9	10	11	12

Test Legend:

1 8270D_S	2 PB_S	3 TPH(D)WSG_S	4	5
6	7	8	9	10
11	12			
The following SampIDs: 001A, 000 testgroup.	2A, 003A, 004A, 005A, 006A, 007A, 008	A, 009A, 010A, 011A, 012A, 013A, 014A, 0	15A contain	Prepared by: Ana Venegas

Comments: SEND HARD COPY

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

Sample Receipt Checklist

Client Name:	Treadwell & Roll	0			Date a	and Time Received:	4/7/2011	7:38:08 PM
Project Name:	#730482302; 5812	2 Hollis St			Check	klist completed and i	eviewed by:	Ana Venegas
WorkOrder N°:	1104220	Matrix Soil			Carrie	er: Rob Pringle (N	IAI Courier)	
		<u>Chair</u>	n of Cu	ıstody (C	COC) Informa	ation_		
Chain of custody	y present?		Yes	V	No 🗆			
Chain of custody	/ signed when relinqui	shed and received?	Yes	V	No \square			
Chain of custody	agrees with sample I	abels?	Yes	✓	No 🗌			
Sample IDs noted	d by Client on COC?		Yes	V	No \square			
Date and Time of	f collection noted by Cli	ient on COC?	Yes	~	No 🗆			
Sampler's name r	noted on COC?		Yes	V	No \square			
		<u>s</u>	ample	Receipt	t Information	<u>1</u>		
Custody seals in	tact on shipping conta	iner/cooler?	Yes		No \square		NA 🔽	
Shipping contain	er/cooler in good cond	lition?	Yes	V	No \square			
Samples in prope	er containers/bottles?		Yes	~	No 🗆			
Sample containe	ers intact?		Yes	✓	No \square			
Sufficient sample	e volume for indicated	test?	Yes	✓	No 🗌			
		Sample Prese	rvatio	n and Ho	old Time (HT) Information		
All samples recei	ived within holding tim	e?	Yes	✓	No 🗌			
Container/Temp I	Blank temperature		Coole	er Temp:	6°C		NA \square	
Water - VOA via	ls have zero headspa	ce / no bubbles?	Yes		No \square	No VOA vials subm	itted 🗹	
Sample labels ch	hecked for correct pre	servation?	Yes	~	No 🗌			
Metal - pH accep	otable upon receipt (pH	l<2)?	Yes		No \square		NA 🗹	
Samples Receive	ed on Ice?		Yes	V	No 🗆			
		(Ісе Тур	e: WE	ET ICE)			
* NOTE: If the "N	No" box is checked, se	ee comments below.						
		======	=					
Client contacted:		Date contac	ted:			Contacted	by:	
Comments:								

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104220

Lab ID	1104220-001A							
Client ID				P5-1				
Matrix				Soil				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
		Surre	gate Re	coveries (%)				
%SS1:	115			%SS2:	100)		
%SS3:	104			%SS4:	78			
					,,			

Comments

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; <math>DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
333 Montgomery St., Suite 1300	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104220

Lab ID				1104220-002A			
Client ID				P5-2			
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surro	gate Re	coveries (%)			
%SS1:	113	3		%SS2:	97		
4					1		

 Surrogate Recoveries (%)

 %SS1:
 113
 %SS2:
 97

 %SS3:
 104
 %SS4:
 80

 %SS5:
 83
 %SS6:
 81

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; <math>DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.

^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
333 Montgomery St., Suite 1300	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/11/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Work Order: 1104220 Extraction Method: SW3550B

			-				
Lab ID				1104220-003A		-	
Client ID				P5-3			
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
			1	·		<u> </u>	1
Acenaphthene	ND ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor Benzidine	ND	1.0	0.33	Anthracene Benzoic Acid	ND	1.0	0.33
	ND	1.0	1.6		ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surre	gate Re	coveries (%)			
%SS1:	92			%SS2:	87		
% SS3·	80			% SSA·	9/		

Surrogate Recoveries (%)									
%SS1:	92	%SS2:	87						
%SS3:	89	% SS4:	94						
%SS5:	76	%SS6:	83						
	·		·						

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS =Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/11/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104220

Lab ID		1104220-004A					
Client ID		P5-4					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33	·			

Surrogate Recoveries (%)					
%SS1:	94	%SS2:	88		
%SS3:	90	%SS4:	96		
%SS5:	82	%SS6:	86		

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS =Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Extraction Method: SW3550B Work Order: 1104220

Lab ID		1104220-005A					
Client ID		P5-5					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33	·			
		Surre	ogate Re	coveries (%)			

Surrogate Recoveries (%)						
%SS1:	92	%SS2:	84			
%SS3:	88	%SS4:	93			
%SS5:	77	%SS6:	82			

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS =Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Extraction Method: SW3550B Work Order: 1104220

Lab ID	1104220-006A						
Client ID				P5-6			
Matrix		Soil					
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surre	gate Re	coveries (%)			
%SS1:	93			%SS2:	85		
							+

95 %SS3: 87 %SS4:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



 $^{*\} water\ samples\ in\ \mu g/k,\ soil/sludge/solid\ samples\ in\ mg/kg,\ wipe\ samples\ in\ \mu g/wipe,\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ TCLP\ \&\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ produ$ SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104220

Lab ID	1104220-007A						
Client ID		P5-7					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1.2.4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
1,2,1 Themoroccuzene							

Surrogate Recoveries (%)					
%SS1:	94	%SS2:	85		
%SS3:	86	%SS4:	96		
%SS5:	75	%SS6:	84		
	· · · · · · · · · · · · · · · · · · ·	·			

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS =Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Extraction Method: SW3550B Work Order: 1104220

Acetochlor ND	Lab ID		1104220-008A					
Compound Concentration DF Reposing Compound Concentration DF Reposing Limit	Client ID				P5-8			
Compound Concentration DF	Matrix				Soil			
Acetochlor ND	Compound	Concentration *	DF		Compound	Concentration *	DF	
Benzidine	Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Benzo(a)anthracene ND	Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzo(k)fluoranthene ND	Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)pyrene ND	Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
1,1-Bipheny ND	Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Bis (2-chloroethyl) Ether ND 1.0 0.33 Bis (2-chloroisopropyl) Ether ND 1.0 0.33 Bis (2-chlylnexyl) Phthalate ND 1.0 0.33 4-Bromophenyl Phenyl Ether ND 1.0 0.33 Bis (2-chlylnexyl) Phthalate ND 1.0 0.33 4-Chloroaniline ND 1.0 0.66 4-Chloro-3-methylphenol ND 1.0 0.33 4-Chloroanphthalene ND 1.0 0.33 2-Chlorophenyl Phenyl Ether ND 1.0 0.33 2-Chlorophenyl Phenyl Ether ND 1.0 0.33 2-Chlorophenol ND 1.0 0.33 3-Chlorophenyl Phenyl Ether ND 1.0 0.33	Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
Bis (2-ethylhexyl) Phthalate	1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Butylbenzyl Phthalate	Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
A-Chloro-3-methylphenol ND 1.0 0.33 2-Chloronaphthalene ND 1.0 0.33 2-Chlorophenol ND 1.0 0.33 4-Chlorophenyl Phenyl Ether ND 1.0 0.33 0.33 1.0 0.33 1.0 0.33 Dibenzo(a,h)anthracene ND 1.0 0.33 0.33 0.33 0.35 0.	Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
2-Chlorophenol ND	Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
Chrysene	4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
ND	2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
1,2-Dichlorobenzene ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.33 1,4-Dichlorobenzene ND 1.0 0.33 3,3-Dichlorobenzidine ND 1.0 0.66 2,4-Dichlorophenol ND 1.0 0.33 Diethyl Phthalate ND 1.0 0.33 2,4-Dimethylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 4,6-Dinitro-2-methylphenol ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0	Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
1,4-Dichlorobenzene ND 1.0 0.33 3,3-Dichlorobenzidine ND 1.0 0.66 2,4-Dichlorophenol ND 1.0 0.33 Diethyl Phthalate ND 1.0 0.33 2,4-Dimethylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 4,6-Dinitro-2-methylphenol ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Fluoranthene ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobenzene ND 1.0	Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
2,4-Dichlorophenol ND 1.0 0.33 Diethyl Phthalate ND 1.0 0.33 2,4-Dimethylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 4,6-Dinitro-2-methylphenol ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 1.6 Hexachlorobutadiene ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylnaphthalene ND 1.0	1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
2,4-Dimethylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 4,6-Dinitro-2-methylphenol ND 1.0 1.6 2,4-Dinitrophenol ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 1.6 Hexachlorocyclopentadiene ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 2-Methylphenol (o-Cresol) ND 1.0 0.33 2-Nitroaniline ND 1.0 </td <td>1,4-Dichlorobenzene</td> <td>ND</td> <td>1.0</td> <td>0.33</td> <td>3,3-Dichlorobenzidine</td> <td>ND</td> <td>1.0</td> <td>0.66</td>	1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
4,6-Dinitro-2-methylphenol ND 1.0 1.6 2,4-Dinitrophenol ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 1.6 Hexachlorobutadiene ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 1.6 Hexachlorocyclopentadiene ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 <	2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 1,0-Diphenylhydrazine ND 1.0 0.33 1,0-Di	2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
Di-n-octyl Phthalate	4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 2-Methylphenol (o-Cresol) ND 1.0 0.33 3 &/or 4-Methylphenol (m,p-Cres ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrosodiphenylamine ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 0.33 Pyrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33 1.0 0.33 ND 1.0 0.33 0.33 1.0 0.33 0.33 0.33 0.33 1.0 0.33 0.33 0.33 0.33 1.0 0.33 0.33 0.33 0.33 1.0 0.33 0.33 0.33 0.33 1.0 0.33 0.33 0.33 1.0 0.33 0.33 0.33 1.0 0.33 0.33 0.33 1.0 0.33 0.33 0.33 1.0 0.33 0.33 0.33 1.0 0.33 0.33 0.33 1.0 0.33 0.33 0.33 1.0 0.33 0.33 0.33 1.0 0.33 0.33 0.33 1.0 0.33 0.33	2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Hexachlorobenzene	Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Hexachlorocyclopentadiene	Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene ND 1.0 0.33 Isophorone ND 1.0 0.33	Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
2-Methylnaphthalene ND 1.0 0.33 2-Methylphenol (o-Cresol) ND 1.0 0.33 3 &/or 4-Methylphenol (m,p-Cres) ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33	Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33	2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33	N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33	Pentachlorophenol	ND	1.0	1.6		ND	1.0	0.33
2.4.6-Trichlorophenol ND 1.0 0.33	Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
	1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
Surrogate Recoveries (%)	2.4.6-Trichlorophenol	ND	1.0	0.33				
			Surre	ogate Re	coveries (%)			

%SS1: 96 %SS2: 85 %SS3: 88 %SS4: 98 %SS6:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



 $^{*\} water\ samples\ in\ \mu g/k,\ soil/sludge/solid\ samples\ in\ mg/kg,\ wipe\ samples\ in\ \mu g/wipe,\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ TCLP\ \&\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ produ$ SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Extraction Method: SW3550B Work Order: 1104220

Lab ID		1104220-009A					
Client ID		P5-9					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surre	ogate Re	coveries (%)			
W 991							

Surrogate Recoveries (%)							
%SS1:	93	%SS2:	84				
%SS3:	85	%SS4:	96				
%SS5:	75	%SS6:	83				

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS =Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Extraction Method: SW3550B Work Order: 1104220

Lab ID		1104220-010A						
Client ID		P5-10						
Matrix		Soil						
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	0.42	1.0	0.33	Benzo(b)fluoranthene	0.60	1.0	0.33	
Benzo(k)fluoranthene	0.49	1.0	0.33	Benzo(g,h,i)perylene	1.2	1.0	0.33	
Benzo(a)pyrene	0.79	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	0.51	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	1.1	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	0.68	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	0.48	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	1.9	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
Surrogate Recoveries (%)								

%SS6:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



¹⁰⁰ %SS1: 97 %SS2 %SS3: 112 %SS4: 89

 $^{*\} water\ samples\ in\ \mu g/k,\ soil/sludge/solid\ samples\ in\ mg/kg,\ wipe\ samples\ in\ \mu g/wipe,\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ TCLP\ \&\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ produ$ SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Work Order: 1104220 Extraction Method: SW3550B

Lab ID		1104220-011A						
Client ID		P5-11						
Matrix				Soil				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND<16	50	0.33	Acenaphthylene	ND<16	50	0.33	
Acetochlor	ND<16	50	0.33	Anthracene	ND<16	50	0.33	
Benzidine	ND<80	50	1.6	Benzoic Acid	ND<80	50	1.6	
Benzo(a)anthracene	ND<16	50	0.33	Benzo(b)fluoranthene	ND<16	50	0.33	
Benzo(k)fluoranthene	ND<16	50	0.33	Benzo(g,h,i)perylene	ND<16	50	0.33	
Benzo(a)pyrene	ND<16	50	0.33	Benzyl Alcohol	ND<80	50	1.6	
1,1-Biphenyl	ND<16	50	0.33	Bis (2-chloroethoxy) Methane	ND<16	50	0.33	
Bis (2-chloroethyl) Ether	ND<16	50	0.33	Bis (2-chloroisopropyl) Ether	ND<16	50	0.33	
Bis (2-ethylhexyl) Phthalate	ND<16	50	0.33	4-Bromophenyl Phenyl Ether	ND<16	50	0.33	
Butylbenzyl Phthalate	ND<16	50	0.33	4-Chloroaniline	ND<33	50	0.66	
4-Chloro-3-methylphenol	ND<16	50	0.33	2-Chloronaphthalene	ND<16	50	0.33	
2-Chlorophenol	ND<16	50	0.33	4-Chlorophenyl Phenyl Ether	ND<16	50	0.33	
Chrysene	ND<16	50	0.33	Dibenzo(a,h)anthracene	ND<16	50	0.33	
Dibenzofuran	ND<16	50	0.33	Di-n-butyl Phthalate	ND<16	50	0.33	
1,2-Dichlorobenzene	ND<16	50	0.33	1,3-Dichlorobenzene	ND<16	50	0.33	
1,4-Dichlorobenzene	ND<16	50	0.33	3,3-Dichlorobenzidine	ND<33	50	0.66	
2,4-Dichlorophenol	ND<16	50	0.33	Diethyl Phthalate	ND<16	50	0.33	
2,4-Dimethylphenol	ND<16	50	0.33	Dimethyl Phthalate	ND<16	50	0.33	
4,6-Dinitro-2-methylphenol	ND<80	50	1.6	2,4-Dinitrophenol	ND<80	50	1.6	
2,4-Dinitrotoluene	ND<16	50	0.33	2,6-Dinitrotoluene	ND<16	50	0.33	
Di-n-octyl Phthalate	ND<16	50	0.33	1,2-Diphenylhydrazine	ND<16	50	0.33	
Fluoranthene	ND<16	50	0.33	Fluorene	ND<16	50	0.33	
Hexachlorobenzene	ND<16	50	0.33	Hexachlorobutadiene	ND<16	50	0.33	
Hexachlorocyclopentadiene	ND<80	50	1.6	Hexachloroethane	ND<16	50	0.33	
Indeno (1,2,3-cd) pyrene	ND<16	50	0.33	Isophorone	ND<16	50	0.33	
2-Methylnaphthalene	ND<16	50	0.33	2-Methylphenol (o-Cresol)	ND<16	50	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND<16	50	0.33	Naphthalene	ND<16	50	0.33	
2-Nitroaniline	ND<80	50	1.6	3-Nitroaniline	ND<80	50	1.6	
4-Nitroaniline	ND<80	50	1.6	Nitrobenzene	ND<16	50	0.33	
2-Nitrophenol	ND<80	50	1.6	4-Nitrophenol	ND<80	50	1.6	
N-Nitrosodiphenylamine	ND<16	50	0.33	N-Nitrosodi-n-propylamine	ND<16	50	0.33	
Pentachlorophenol	ND<80	50	1.6	Phenanthrene	ND<16	50	0.33	
Phenol	ND<16	50	0.33	Pyrene	ND<16	50	0.33	
1,2,4-Trichlorobenzene	ND<16	50	0.33	2,4,5-Trichlorophenol	ND<16	50	0.33	
2.4.6-Trichlorophenol	ND<16	50	0.33					
		Surre	gate Re	coveries (%)				

%SS1: 59 %SS2: %SS3: 56 %SS4: 92

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



 $^{*\} water\ samples\ in\ \mu g/k,\ soil/sludge/solid\ samples\ in\ mg/kg,\ wipe\ samples\ in\ \mu g/wipe,\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ TCLP\ \&\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ produ$ SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Extraction Method: SW3550B Work Order: 1104220

Lab ID		1104220-012A						
Client ID		P5-12						
Matrix				Soil				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND 1.0 0.33							
	Surrogate Recoveries (%)							
%SS1:	94			%SS2:	84			

%SS3: 85 %SS4 96

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



 $^{*\} water\ samples\ in\ \mu g/k,\ soil/sludge/solid\ samples\ in\ mg/kg,\ wipe\ samples\ in\ \mu g/wipe,\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ TCLP\ \&\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ produ$ SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/09/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Extraction Method: SW3550B Work Order: 1104220

Lab ID		1104220-013A						
Client ID		P5-13						
Matrix				Soil				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
		Surro	gate Re	coveries (%)				
%SS1:	97			%SS2:	97			
21 000				21.001	-			

%SS3

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



 $^{*\} water\ samples\ in\ \mu g/k,\ soil/sludge/solid\ samples\ in\ mg/kg,\ wipe\ samples\ in\ \mu g/wipe,\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ TCLP\ \&\ product/oil/non-aqueous\ liquid\ samples\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ product/oil/non-aqueous\ and\ all\ product/oil/non-aqueous\ and\ and\ produ$ SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Work Order: 1104220 Extraction Method: SW3550B

Lab ID	1104220-014A							
Client ID		P5-14						
Matrix				Soil				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
		Surre	ogate Re	coveries (%)				
%SS1:	96			%SS2:	84			

%SS1: 96 %SS2: 84 %SS3: 89 %SS4: 95 %SS5: 95 95	Surrogate Recoveries (%)							
	%SS1:	96	%SS2:	84				
V 995	%SS3:	89	%SS4:	95				
<u> </u>	%SS5:	81	%SS6:	85				

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS =Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/12/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104220

Lab ID		1104220-015A						
Client ID		P5-15						
Matrix		Soil						
Compound	Concentration *	DF	Reporting Limit	Compound Concentration		DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND ND	1.0	1.6	3-Nitroaniline	ND ND	1.0	1.6	
4-Nitroaniline	ND ND	1.0	1.6	Nitrobenzene	ND ND	1.0	0.33	
2-Nitrophenol	ND ND	1.0	1.6	4-Nitrophenol	ND ND	1.0	1.6	
N-Nitrosodiphenylamine	ND ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND ND	1.0	0.33	
Pentachlorophenol	ND ND	1.0	1.6	Phenanthrene	ND ND	1.0	0.33	
Phenol	ND ND	1.0	0.33	Pyrene	ND ND	1.0	0.33	
		1.0		·				
1,2,4-Trichlorobenzene	ND ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND			coveries (%)				

Surrogate Recoveries (%)						
%SS1:	94	%SS2:	80			
%SS3:	86	%SS4:	96			
%SS5:	78	%SS6:	82			

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS =Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled: 04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/07/11
	Client Contact: Peter Cusack	Date Extracted: 04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 04/08/11-04/12/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline*

Extraction method SW5	030B	Analytical metho	ds SW8015Bm	Wo	ork Order:	1104220	
Lab ID	Client ID	Matrix	TPH(g)	DF	% SS	Comment	
001A	P5-1	S	ND	1	84		
002A	P5-2	S	ND	1	85		
003A	P5-3	s	ND	1	78		
004A	P5-4	S	ND	1	80		
005A	P5-5	s	ND	1	87		
006A	P5-6	S	ND	1	95		
007A	P5-7	S	ND	1	94		
008A	P5-8	s	ND	1	93		
009A	P5-9	s	ND	1	81		
010A	P5-10	s	ND	1	102		
011A	P5-11	S	580	20	96	d7,d9	
012A	P5-12	S	ND	1	86		
013A	P5-13	s	ND	1	80		
014A	P5-14	S	ND	1	87		
015A	P5-15	S	ND	1	86		
	ng Limit for DF =1;	W	NA		NA		
	ins not detected at or	S	1.0		mg/K	g	

ND means not detected at or above the reporting limit	S	1.0	mg/Kg
* water and vapor samples are reported in ug/L, soil/sluc	lge/solid sampl	es in mg/kg, wipe samples in ug/wipe, product	/oil/non-aqueous liquid

^{*} water and vapor samples are reported in µg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts in mg/L.

d7) strongly aged gasoline or diesel range compounds are significant in the TPH(g) chromatogram d9) no recognizable pattern

Angela Rydelius, Lab Manager

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:



Extraction method: SW3050B

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled:	04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received:	04/07/11
	Client Contact: Peter Cusack	Date Extracted:	04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/08/11-04/11/11

Lead by ICP*

Analytical methods: SW6010B

Lab ID Client ID Matrix Extraction Type DF % SS Comments 1104220-001A P5-1 S TOTAL 105 7.8 1 1104220-002A P5-2 S TOTAL 8.6 1 106 1104220-003A S TOTAL P5-3 ND 1 94 1104220-004A S TOTAL P5-4 7.8 91 1 1104220-005A P5-5 S TOTAL 8.9 104 1104220-006A P5-6 S TOTAL 10 106 1104220-007A P5-7 S TOTAL 5.1 102 1 S 1104220-008A P5-8 TOTAL 14 1 100 1104220-009A P5-9 S TOTAL 8.6 1 102 1104220-010A P5-10 S TOTAL 15 1 103 1104220-011A P5-11 S TOTAL 7.5 1 99 \mathbf{S} 1104220-012A P5-12 TOTAL 8.3 1 100

Reporting Limit for DF =1;	W	TOTAL	NA	μg/L
ND means not detected at or above the reporting limit	S	TOTAL	5.0	mg/Kg

TOTAL

TOTAL

6.4

6.2

*water samples are reported in μ g/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, filter samples in μ g/filter.

 \mathbf{S}

S

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of $0.45 \mu m$ filtered and acidified sample.

P5-13

P5-14

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

1104220-013A

1104220-014A

Angela Rydelius, Lab Manager

94

99

1

1

DHS ELAP Certification 1644

Work Order: 1104220



"When Quality Counts"			16	iepnone: a	8//-252-9262 Fax: 925	-232-9209			
Treadwell & F	Rollo	Client Project Hollis St	ct ID: #	730482302; 5812	2	Date Sampled:	04/05/11		
555 Montgomery St., Suite 1300		Homs St	Homs St			Date Received:	04/07/11		
	3	Client Cont	tact: Per	ter Cusack		Date Extracted:	04/07/11		
San Francisco	o, CA 94111	Client P.O.:				Date Analyzed:	04/08/11	-04/11/1	1
			Le	ad by ICP*					
Extraction method:	: SW3050B		Analy	tical methods: SW60)10B			Work Ord	ler: 1104220
Lab ID	Client ID		Matrix	Extraction Type		Lead	DF	% SS	Comments
1104220-015A	P5-15		S	TOTAL		8.4	1	106	

Reporting Limit for DF =1;	W	TOTAL	NA	μg/L
ND means not detected at or above the reporting limit	S	TOTAL	5.0	mg/Kg

*water samples are reported in μ g/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, filter samples in μ g/filter.

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of $0.45 \mu m$ filtered and acidified sample.

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

Angela Rydelius, Lab Manager

DHS ELAP Certification 1644



Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled:	04/05/11
555 Montgomery St., Suite 1300	Hollis St	Date Received:	04/07/11
	Client Contact: Peter Cusack	Date Extracted:	04/07/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/08/11-04/14/11

Total Extractable Petroleum Hydrocarbons with Silica Gel Clean-Up*

Extraction method: SW3	550B/3630C	Analytical	l methods: SW8015B		Work O			
Lab ID	Client ID	Matrix	TPH-Diesel (C10-C23)	TPH-Motor Oil (C18-C36)	DF	% SS	Comments	
1104220-001A	P5-1	S	ND	ND	1	103		
1104220-002A	P5-2	S	ND	ND	1	106		
1104220-003A	P5-3	S	ND	ND	1	104		
1104220-004A	P5-4	S	ND	ND	1	112		
1104220-005A	P5-5	S	ND	ND	1	105		
1104220-006A	P5-6	S	ND	ND	1	104		
1104220-007A	P5-7	S	ND	ND	1	118		
1104220-008A	P5-8	S	1.6	6.9	1	106	e7,e2	
1104220-009A	P5-9	S	6.5	14	1	105	e7,e2	
1104220-010A	P5-10	S	8.9	11	1	107	e7,e2	
1104220-011A	P5-11	S	13,000	5500	100	104	e1	
1104220-012A	P5-12	S	4.0	ND	1	105	e1	
1104220-013A	P5-13	S	2.4	ND	1	104	e2	
1104220-014A	P5-14	S	ND	ND	1	106		
1104220-015A	P5-15	S	ND	ND	1	106		
Reportin	g Limit for DF =1;	W	NA	NA		ug/.	L	

ND manns not detected at an		1171	1171	ugL		
ND means not detected at or	C	1.0	5.0	mg/Kg		
above the reporting limit	_ S	1.0	3.0	nig/Kg		
		, ,,,,,				

^{*} water samples are reported in μ g/L, wipe samples in μ g/wipe, soil/solid/sludge samples in mg/kg, product/oil/non-aqueous liquid samples in mg/L, and all DISTLC / STLC / SPLP / TCLP extracts are reported in μ g/L.

%SS = Percent Recovery of Surrogate Standard. DF = Dilution Factor

- e1) unmodified or weakly modified diesel is significant
- e2) diesel range compounds are significant; no recognizable pattern
- e7) oil range compounds are significant



[#] cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract.

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

QC SUMMARY REPORT FOR SW8270C

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57506 WorkOrder 1104220

EPA Method SW8270C Extraction SW3550B Spiked Sample ID: 1104143-001A)01A			
Analyte	Sample	Spiked	MS	MSD	MSD MS-MSD LCS LCSD LCS-LCSD Acceptance Cri					Criteria (%)		
7 may to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Acenaphthene	ND	2	61.2	64.9	5.81	64.8	65.1	0.431	30 - 130	30	30 - 130	30
4-Chloro-3-methylphenol	ND	4	59.8	63.7	6.32	72	70.3	2.34	30 - 130	30	30 - 130	30
2-Chlorophenol	ND	4	60.8	57.1	6.31	61.7	62.9	1.97	30 - 130	30	30 - 130	30
1,4-Dichlorobenzene	ND	2	66.4	70.3	5.76	68.3	67.9	0.543	30 - 130	30	30 - 130	30
2,4-Dinitrotoluene	ND	2	61.2	68.1	10.7	74.3	72.5	2.44	30 - 130	30	30 - 130	30
4-Nitrophenol	ND	4	89.2	89.6	0.492	72.1	67.9	5.98	30 - 130	30	30 - 130	30
N-Nitrosodi-n-propylamine	ND	2	59.8	53.4	11.4	55.6	54.8	1.47	30 - 130	30	30 - 130	30
Pentachlorophenol	ND	4	54	53.9	0.297	46.6	45.6	2.30	30 - 130	30	30 - 130	30
Phenol	ND	4	61.4	59.8	2.65	63.3	62	2.02	30 - 130	30	30 - 130	30
Pyrene	ND	2	49.7	50.2	1.00	62.5	54.6	13.6	30 - 130	30	30 - 130	30
1,2,4-Trichlorobenzene	ND	2	74.7	73.9	1.02	79.3	77.3	2.53	30 - 130	30	30 - 130	30
%SS1:	81	200	76	73	4.02	84	83	1.80	30 - 130	30	30 - 130	30
%SS2:	77	200	72	73	0.730	89	87	1.85	30 - 130	30	30 - 130	30
%SS3:	82	200	83	79	4.30	96	94	2.41	30 - 130	30	30 - 130	30
% SS4:	80	200	80	80	0	88	89	0.587	30 - 130	30	30 - 130	30
%SS5:	67	200	71	66	6.85	98	95	3.22	30 - 130	30	30 - 130	30
%SS6:	74	200	74	78	5.72	104	94	9.34	30 - 130	30	30 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57506 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104220-001A	04/05/11 9:30 AM	04/07/11	04/12/11 12:41 PM	1104220-002A	04/05/11 1:21 PM	04/07/11	04/12/11 2:04 PM
1104220-003A	04/05/11 1:24 PM	04/07/11	04/11/11 10:05 PM	1104220-004A	04/05/11 1:30 PM	04/07/11	04/11/11 11:22 PM
1104220-005A	04/05/11 1:35 PM	04/07/11	04/12/11 12:38 AM	1104220-006A	04/05/11 10:30 AM	04/07/11	04/12/11 1:54 AM
1104220-007A	04/05/11 10:40 AM	04/07/11	04/12/11 3:09 AM	1104220-008A	04/05/11 10:50 AM	04/07/11	04/12/11 4:24 AM
1104220-009A	04/05/11 9:25 AM	04/07/11	04/12/11 5:38 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and / or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix, sample diluted due to high matrix or analyte content, or MS/MSD samples diluted due to high organic content.

#) surrogate diluted out of range; & = low or no recovery of surrogate or target analytes due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.

QC SUMMARY REPORT FOR SW8270C

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57564 WorkOrder 1104220

EPA Method SW8270C	Extra	ction SW	3550B					S	Spiked Sar	nple ID	1104220-0	13A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acc	eptance	Criteria (%)	1
Analyto	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Acenaphthene	ND	2	73.6	75.7	2.83	75.6	75.2	0.451	30 - 130	30	30 - 130	30
4-Chloro-3-methylphenol	ND	4	85.9	88.3	2.82	75.2	79.2	5.27	30 - 130	30	30 - 130	30
2-Chlorophenol	ND	4	73.5	75.3	2.39	73.8	75.2	1.87	30 - 130	30	30 - 130	30
1,4-Dichlorobenzene	ND	2	75.2	77.7	3.22	76.4	78.7	3.03	30 - 130	30	30 - 130	30
2,4-Dinitrotoluene	ND	2	79.8	82	2.68	77.4	81.6	5.29	30 - 130	30	30 - 130	30
4-Nitrophenol	ND	4	87.5	87.8	0.371	71.3	71.1	0.309	30 - 130	30	30 - 130	30
N-Nitrosodi-n-propylamine	ND	2	88.1	92.8	5.17	85.3	88.3	3.43	30 - 130	30	30 - 130	30
Pentachlorophenol	ND	4	60.6	63.2	4.26	32.5	30.5	6.30	30 - 130	30	30 - 130	30
Phenol	ND	4	70.4	72.8	3.28	67.8	69.4	2.27	30 - 130	30	30 - 130	30
Pyrene	ND	2	75.2	77.2	2.63	72.3	77.7	7.24	30 - 130	30	30 - 130	30
1,2,4-Trichlorobenzene	ND	2	85.8	88	2.57	79.8	87.9	9.71	30 - 130	30	30 - 130	30
%SS1:	97	200	93	95	2.04	92	93	1.86	30 - 130	30	30 - 130	30
%SS2:	97	200	95	95	0	93	95	2.70	30 - 130	30	30 - 130	30
%SS3:	99	200	101	103	1.63	98	104	6.67	30 - 130	30	30 - 130	30
%SS4:	88	200	88	90	2.77	97	91	6.36	30 - 130	30	30 - 130	30
%SS5:	102	200	96	101	5.66	82	89	7.50	30 - 130	30	30 - 130	30
%SS6:	91	200	90	91	1.75	88	93	6.07	30 - 130	30	30 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57564 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104220-010A	04/05/11 9:20 AM	04/07/11	04/09/11 3:54 AM	1104220-011A	04/05/11 9:15 AM	04/07/11	04/12/11 7:39 PM
1104220-012A	04/05/11 9:10 AM	04/07/11	04/12/11 6:52 AM	1104220-013A	04/05/11 9:05 AM	04/07/11	04/09/11 12:25 PM
1104220-014A	04/05/11 1:46 PM	04/07/11	04/12/11 8:07 AM	1104220-015A	04/05/11 9:40 AM	04/07/11	04/12/11 9:23 AM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and / or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix, sample diluted due to high matrix or analyte content, or MS/MSD samples diluted due to high organic content.

#) surrogate diluted out of range; & = low or no recovery of surrogate or target analytes due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.

QC SUMMARY REPORT FOR SW8021B/8015Bm

QC Matrix: Soil BatchID: 57513 WorkOrder 1104220 W.O. Sample Matrix: Soil

EPA Method SW8015Bm	Extra	ction SW	5030B					5	Spiked San	nple ID:	: 1104152-0	03A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
Analyte	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex [£]	ND	0.60	124	125	0.352	128	124	2.79	70 - 130	20	70 - 130	20
MTBE	ND	0.10	101	109	7.37	108	112	3.44	70 - 130	20	70 - 130	20
Benzene	ND	0.10	89.5	90.4	1.03	91.1	93.6	2.67	70 - 130	20	70 - 130	20
Toluene	ND	0.10	88.1	88.8	0.737	88.7	91.8	3.39	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	89.5	90.6	1.24	90.8	94.3	3.81	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	89.5	90.9	1.51	90.6	93.9	3.57	70 - 130	20	70 - 130	20
%SS:	83	0.10	78	82	4.96	76	87	12.8	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57513 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104220-001A	04/05/11 9:30 AM	04/07/11	04/09/11 2:30 AM	1104220-002A	04/05/11 1:21 PM	04/07/11	04/09/11 2:59 AM
1104220-003A	04/05/11 1:24 PM	04/07/11	04/09/11 4:27 AM	1104220-004A	04/05/11 1:30 PM	04/07/11	04/09/11 4:57 AM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QC SUMMARY REPORT FOR SW8021B/8015Bm

QC Matrix: Soil BatchID: 57563 WorkOrder 1104220 W.O. Sample Matrix: Soil

EPA Method SW8015Bm	Extra	ction SW	5030B					5	Spiked San	nple ID:	: 1104220-0)14A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%))
7 tildiy to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex ^f)	ND	0.60	120	113	5.98	120	121	1.02	70 - 130	20	70 - 130	20
MTBE	ND	0.10	92.1	94.8	2.85	94.5	87.4	7.84	70 - 130	20	70 - 130	20
Benzene	ND	0.10	89.9	90.6	0.773	93.4	89.5	4.25	70 - 130	20	70 - 130	20
Toluene	ND	0.10	88.1	89.5	1.60	90.4	87.5	3.29	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	89.3	91.5	2.46	91.5	88.5	3.42	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	89.4	91.3	2.07	91.1	88.2	3.25	70 - 130	20	70 - 130	20
%SS:	87	0.10	88	94	6.43	82	84	2.45	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57563 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104220-005A	04/05/11 1:35 PM	04/07/11	04/09/11 5:26 AM	1104220-006A	04/05/11 10:30 AM	04/07/11	04/08/11 5:36 PM
1104220-007A	04/05/11 10:40 AM	04/07/11	04/08/11 6:36 PM	1104220-008A	04/05/11 10:50 AM	04/07/11	04/08/11 7:07 PM
1104220-009A	04/05/11 9:25 AM	04/07/11	04/09/11 7:23 AM	1104220-010A	04/05/11 9:20 AM	04/07/11	04/12/11 12:35 AM
1104220-011A	04/05/11 9:15 AM	04/07/11	04/08/11 11:33 PM	1104220-012A	04/05/11 9:10 AM	04/07/11	04/12/11 12:48 AM
1104220-013A	04/05/11 9:05 AM	04/07/11	04/12/11 3:15 AM	1104220-014A	04/05/11 1:46 PM	04/07/11	04/12/11 3:44 AM
1104220-015A	04/05/11 9:40 AM	04/07/11	04/09/11 8:22 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104220

EPA Method SW6010B		Extraction SW3050B					BatchID	: 57460	Spiked Sample ID: 1104093-003A							
Analyte	Sample	Spiked	MS	MSD	MS-MSD	Spiked	LCS	LCSD	LCS-LCSD	Acce	eptance	e Criteria (%	o)			
7 illuly to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD			
Lead	15	50	96.9	104	5.61	10	101	95.6	5.17	75 - 125	25	75 - 125	25			
%SS:	98	500	98	95	2.64	500	103	100	3.84	70 - 130	20	70 - 130	20			

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57460 SUMMARY

Lab ID	Date Sampled	Date Extracte	d Date Analyzed	Lab ID	Date Sampled	Date Extracte	d Date Analyzed
1104220-001A	04/05/11 9:30 AM	04/07/11	04/11/11 11:09 AM	1104220-002A	04/05/11 1:21 PM	04/07/11	04/11/11 11:11 AM
1104220-003A	04/05/11 1:24 PM	04/07/11	04/08/11 3:12 PM	1104220-004A	04/05/11 1:30 PM	04/07/11	04/08/11 3:14 PM
1104220-005A	04/05/11 1:35 PM	04/07/11	04/11/11 11:13 AM	1104220-006A	04/05/11 10:30 AM	04/07/11	04/11/11 11:16 AM
1104220-007A	04/05/11 10:40 AM	04/07/11	04/08/11 3:16 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104220

EPA Method SW6010B			Extract	ion SW3	3050B		BatchID	: 57565	Spiked Sample ID: 1104220-012A						
Analyte	Sample	Spiked	MS	MSD	MS-MSD	Spiked	LCS	LCSD	LCS-LCSD	Acce	eptance	otance Criteria (%)			
7 mary to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD		
Lead	8.3	50	83.3	86.1	2.77	10	85.8	87.8	2.27	75 - 125	25	75 - 125	25		
%SS:	100	500 98 102 4.03					96	98	1.95	70 - 130	20	70 - 130	20		

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57565 SUMMARY

Lab ID	Date Sampled	Date Extracted	d Date Analyzed	Lab ID	Date Sampled	Date Extract	ed Date Analyzed
1104220-008A	04/05/11 10:50 AM	04/07/11	04/11/11 11:18 AM	1104220-009A	04/05/11 9:25 AM	04/07/11	04/11/11 11:20 AM
1104220-010A	04/05/11 9:20 AM	04/07/11	04/11/11 11:23 AM	1104220-011A	04/05/11 9:15 AM	04/07/11	04/08/11 3:19 PM
1104220-012A	04/05/11 9:10 AM	04/07/11	04/08/11 2:42 PM	1104220-013A	04/05/11 9:05 AM	04/07/11	04/08/11 3:21 PM
1104220-014A	04/05/11 1:46 PM	04/07/11	04/08/11 3:23 PM	1104220-015A	04/05/11 9:40 AM	04/07/11	04/11/11 11:25 AM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57512 WorkOrder 1104220

EPA Method SW8015B	Extrac	Spiked Sample ID: 1104154-002A										
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
, and y to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	2.5	40	119	118	1.02	115	113	2.17	70 - 130	30	70 - 130	30
%SS:	106	25	108	107	0.632	95	93	2.07	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57512 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104220-001A	04/05/11 9:30 AM	04/07/11	04/09/11 1:31 AM	1104220-002A	04/05/11 1:21 PM	04/07/11	04/11/11 9:08 PM
1104220-003A	04/05/11 1:24 PM	04/07/11	04/09/11 4:51 AM	1104220-004A	04/05/11 1:30 PM	04/07/11	04/08/11 8:51 AM
1104220-005A	04/05/11 1:35 PM	04/07/11	04/09/11 2:38 AM	1104220-006A	04/05/11 10:30 AM	04/07/11	04/09/11 3:45 AM
1104220-007A	04/05/11 10:40 AM	04/07/11	04/13/11 12:46 AM	1104220-008A	04/05/11 10:50 AM	04/07/11	04/14/11 1:45 PM
1104220-009A	04/05/11 9:25 AM	04/07/11	04/13/11 1:20 AM	1104220-010A	04/05/11 9:20 AM	04/07/11	04/08/11 8:59 PM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

DHS ELAP Certification 1644

QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57562 WorkOrder 1104220

EPA Method SW8015B	Extrac		Spiked Sample ID: 1104220-015A										
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)		
, and y to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD	
TPH-Diesel (C10-C23)	ND	40	117	116	0.711	118	117	0.817	70 - 130	30	70 - 130	30	
%SS:	106	25	109	109	0	101	101	0	70 - 130	30	70 - 130	30	

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57562 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104220-011A	04/05/11 9:15 AM	04/07/11	04/12/11 1:14 AM	1104220-012A	04/05/11 9:10 AM	04/07/11	04/08/11 3:38 PM
1104220-013A	04/05/11 9:05 AM	04/07/11	04/08/11 2:30 PM	1104220-014A	04/05/11 1:46 PM	04/07/11	04/12/11 5:59 PM
1104220-015A	04/05/11 9:40 AM	04/07/11	04/12/11 7:08 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

DHS ELAP Certification 1644

McCampbell Analytical,	Inc.
"When Quality Counts"	

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300		Date Received: 04/15/11
oss frongomery bu, butter 1900	Client Contact: Peter Cusack	Date Reported: 04/19/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/19/11

WorkOrder: 1104453

April 20, 2011

D	D .
I lear	Peter:

Enclosed within are:

- 1) The results of the 9 analyzed samples from your project: #730482302; 5812 Hollis St.,
- 2) A QC report for the above samples,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.



PRESERVATION

CHAIN OF CUSTODY RECORD

Environmental and Geote	echnical Consu	ltant								00, Sa											5.9	55.9041		× .	3	
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Job Number:	130 4	B2305	^							*		70		A	nalys	is R	equ	ieste	d					Turn	around	ı
Project Manager\Cor Samplers:	Kan	Veler Vilani	Sugack								-M-	8												72	ime	115
Recorder (Signature	Required):	-	un				No	. Co	ntai	ners	b	1								1 2				MARIE CO.	1,0	
		•	75.55	N	latri	х	&	Pres	erv	ative	0	1	7	1						clean-up	5					
Field Sample Identification No.	Date	Time	Lab Sample No.	Soil	Water	Other	HCL HCL	HNOS	lce	Other	3		2							Silica del	5	РІОН	F	Remarks		
B-1-15	4-15-11	1350		X		\top		Т	X		X	k	\backslash	П		П	\top			A						
B-2-15	1	13.55		7		\top	\top	+	X		X	X	×	\Box	\top	\forall	+		\vdash	X						
12-3-15		HOD.		×	~	+	+	+	X		V	V	2	\forall	+	\vdash	+	+	H	8	+	+				
13-4-15		1403		X		\top	+	+	2		k	X	K		+	\forall	+	+	\forall	\rightarrow	+					
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ICE/t° CONDIT GOOD CONDIT HEAD SPACE A DECHLORINA	TION ABSENT TED IN LAB	White Copy - APPROPRI CONTAINE PRESERVI	ERSED IN LAB		Yello	ow C	ору	- La	bora	atory				F	ink C	ору -	Fie	ld		(CC	OC Number:	00)453	37	
	VUAS U	CATTER CATE																								

McCampbell Analytical, Inc.

1534 Willow Pass Rd Pittsburg, CA 94565-1701

CHAIN-OF-CUSTODY RECORD

ClientCode: TWRF

WorkOrder: 1104453

Page 1 of 1

(723) 232-7202	WaterTrax	WriteOn	☐ EDF	Excel	Fax	✓ Email	HardCopy	ThirdParty	☐ J-flag
Report to:				Bil	I to:		Req	uested TAT:	3 days
Peter Cusack Treadwell & Rollo 555 Montgomery St., Suite 1300 San Francisco, CA 94111 (415) 955-5244 FAX (415) 955-904	cc: PO: ProjectNo: #	ojcusack@treadv #730482302; 58 ²			•	•	50	te Received: te Printed:	04/15/2011 04/15/2011

				Ī	Requested Tests (See legend below)											
Lab ID	Client ID	Matrix	Collection Date F	lold	1	2	3	4	5	6	7	8	9	10	11	12
1104453-001	B-1-15	Soil	4/15/2011 13:50	П	Α	Α	Α									
1104453-002	B-2-15	Soil	4/15/2011 13:55		Α	Α	Α									
1104453-003	B-3-15	Soil	4/15/2011 14:00		Α	Α	Α									
1104453-004	B-4-15	Soil	4/15/2011 14:05		Α	Α	Α									
1104453-005	B-5-15	Soil	4/15/2011 14:10		Α	Α	Α									
1104453-006	B-6-15	Soil	4/15/2011 14:15		Α	Α	Α									
1104453-007	B-7-15	Soil	4/15/2011 14:20		Α	Α	Α									
1104453-008	B-8-15	Soil	4/15/2011 14:25		Α	Α	Α									
1104453-009	B-9-15	Soil	4/15/2011 14:30		Α	Α	Α									

Test Legend:

1 8270D_S	2 PB_S	3 TPH(DMO)WSG_S	4	5					
6	7	8	9	10					
11	12								
The following SampIDs: 001A, 002A, 003A, 004A, 005A, 006A, 007A, 008A, 009A contain testgroup. Prepared by: Maria Venego									

Comments: SEND HARD COPY. 72hr Rush

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

Sample Receipt Checklist

Client Name: Treadwell & Rollo						Date	and Time Received:	4/15/2011	5:43:36 PM
Project Name:	#730482302;	5812 Hollis	St.			Chec	klist completed and	reviewed by:	Maria Venegas
WorkOrder N°:	1104453	Matrix S	<u>Soil</u>			Carrie	er: Rob Pringle (N	MAI Courier)	
			Chain	of Cu	stody (C	COC) Inform	<u>ation</u>		
Chain of custod	ly present?			Yes	V	No 🗆			
Chain of custod	y signed when re	linquished and	received?	Yes	V	No 🗆			
Chain of custod	ly agrees with sar	nple labels?		Yes	✓	No 🗌			
Sample IDs note	ed by Client on CO	C?		Yes	V	No 🗆			
Date and Time of	of collection noted	by Client on CO	C?	Yes	V	No 🗆			
Sampler's name	noted on COC?			Yes	V	No 🗆			
			Sa	ample	Receipt	t Informatio	<u>n</u>		
Custody seals in	ntact on shipping	container/coole		Yes		No 🗆		NA 🗹	
Shipping contain	ner/cooler in good	condition?		Yes	V	No 🗆			
Samples in prop	per containers/bot	tles?		Yes	V	No 🗆			
Sample contain	ers intact?			Yes	✓	No 🗆			
Sufficient sampl	le volume for indic	cated test?		Yes	✓	No 🗌			
		Sam	ple Preser	vatio	n and Ho	old Time (HT	「) Information		
All samples rece	eived within holdin	ng time?		Yes	✓	No 🗌			
Container/Temp	Blank temperatur	e		Coole	er Temp:	6.8°C		NA 🗆	
Water - VOA via	als have zero hea	dspace / no bul	bbles?	Yes		No 🗆	No VOA vials subr	nitted 🗹	
Sample labels of	checked for correc	ct preservation?		Yes	~	No 🗌			
Metal - pH acce	ptable upon receip	ot (pH<2)?		Yes		No 🗆		NA 🗹	
Samples Receiv	ved on Ice?			Yes	V	No 🗆			
			(Ice Type	e: WE	TICE)			
* NOTE: If the "	'No" box is check	ed, see comme	nts below.						
	=====	=====					=====		
Client contacted: Date contact		ed:			Contacted	d by:			
Comments:									

McCampbell Analytical, Inc.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
555 Workgomery St., State 1500	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/19/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104453

Lab ID	1104453-001A								
Client ID									
Matrix				Soil					
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit		
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33		
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33		
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6		
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33		
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33		
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6		
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33		
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33		
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33		
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66		
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33		
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33		
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33		
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33		
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33		
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66		
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33		
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33		
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6		
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33		
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33		
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33		
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33		
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33		
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33		
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33		
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33		
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6		
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33		
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6		
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33		
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33		
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33		
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33		
2.4.6-Trichlorophenol	ND	1.0	0.33						
		Surro	gate Re	coveries (%)	-				
%SS1:	87			%SS2:	83				
	07				0.5				

Surrogate Recoveries (%) %SS1: 87 %SS2: 83 %SS3: 93 %SS4: 84 %SS5: 78 %SS6: 92

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.

^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/18/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104453

Lab ID		1104453-002A						
Client ID				B-2-15				
Matrix				Soil				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
		Surre	ogate Re	coveries (%)				
<u> </u>			-		1			

Surrogate Recoveries (%)						
%SS1:	96	%SS2:	102			
%SS3:	94	%SS4:	79			
%SS5:	96	%SS6:	94			

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

^{*} water samples in $\mu g/L$, soil/sludge/solid samples in mg/kg, wipe samples in $\mu g/wipe$, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

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Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
333 Wongomery St., Suite 1300	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/18/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104453

Lab ID		1104453-003A						
Client ID				B-3-15				
Matrix				Soil				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
		Surre	gate Re	coveries (%)				
·			TO		T			

Surrogate Recoveries (%)					
%SS1:	96	%SS2:	96		
%SS3:	97	%SS4:	79		
%SS5:	100	%SS6:	98		

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

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The state of the s	1	
Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/15/11
555 Montgomery St. Suite 1300	Hollis St.	Date Received: 04/15/11
555 Montgomery St., Suite 1300	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/18/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104453

Lab ID		1104453-004A					
Client ID				B-4-15			
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surro	gate Re	coveries (%)			
%SS1:	93			%SS2:	91		

%SS1: 93 %SS2: 91 %SS3: 94 %SS4: 82 %SS5: 93 %SS6: 98

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

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Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/18/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104453

Lab ID		1104453-005A					
Client ID				B-5-15			
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				-
		Surro	gate Re	coveries (%)			
%SS1:	96			%SS2:	97		

%SS5:

%SS3:

%SS4:

96

96

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.

82

^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/18/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104453

Lab ID		1104453-006A					
Client ID		B-6-15					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surre	gate Re	coveries (%)			
%SS1:	91			%SS2:	91		
%SS3:	94			%SS4:	83		

Comments

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
333 Wongomery St., State 1300	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/19/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104453

Lab ID 1104453-007A							
Client ID				B-7-15			
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33		<u> </u>		
				coveries (%)			

Surrogate Recoveries (%)					
%SS1:	91	%SS2:	92		
%SS3:	93	%SS4:	83		
%SS5:	94	%SS6:	90		

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	•	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
555 Workgomery St., State 1500	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/19/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104453

Lab ID 1104453-008A							
Client ID		B-8-15					
Matrix		Soil					
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surro	gate Re	coveries (%)			
% SS1·							

Surrogate Recoveries (%)					
%SS1:	90	%SS2:	92		
%SS3:	97	%SS4:	80		
%SS5:	86	%SS6:	94		

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
333 Montgomery St., State 1300	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/19/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104453

Lab ID 1104453-009A							
Client ID				B-9-15			
Matrix		Soil					
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surre	gate Re	coveries (%)			
w.cc1.							

Surrogate Recoveries (%)					
%SS1:	90	%SS2:	90		
%SS3:	92	%SS4:	79		
%SS5:	95	%SS6:	97		

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in $\mu g/L$, soil/sludge/solid samples in mg/kg, wipe samples in $\mu g/wipe$, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
<i>5</i> ,	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 04/15/11-04/19/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline* Analytical methods SW8015Bm Extraction method SW5030B Work Order: 1104453 Lab ID Client ID Matrix TPH(g)DF % SS Comments 001A B-1-15 S 5.2 002A S B-2-15 ND 94 003A S 98 B-3-15 ND 1 004A B-4-15 S ND 1 90 005A B-5-15 S ND 1 100 006A B-6-15 S ND 1 99 007A B-7-15 S 4.1 1 84 d7 008A B-8-15 S ND 1 96 009A B-9-15 S ND 1 Reporting Limit for DF =1; NA

above the reporting limit	S	1.0	mg/Kg		
* water and vapor samples are reported in µg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous liquid					
samples and all TCLP & SPLP extracts in mg/L.					

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

d7) strongly aged gasoline or diesel range compounds are significant in the TPH(g) chromatogram

Angela Rydelius, Lab Manager

ND means not detected at or

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:



Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/15/11
•	Client Contact: Peter Cusack	Date Extracted: 04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/18/11

Lead by ICP*

Extraction method: SW3050B Analytical methods: SW6010B Work Order: 1104453 Lab ID Client ID Matrix Extraction Type Lead DF % SS Comments 1104453-001A S TOTAL 107 B-1-15 6.6 1 1104453-002A B-2-15 S TOTAL 5.6 1 108 1104453-003A S TOTAL B-3-15 8.4 1 107 1104453-004A S TOTAL B-4-15 12 1 108 1104453-005A B-5-15 S TOTAL 108 1104453-006A B-6-15 S TOTAL 105 5.7 1104453-007A B-7-15 S TOTAL 1 108 8.6 \mathbf{S} 1104453-008A TOTAL 1 108 B-8-15 11 1104453-009A B-9-15 S TOTAL 7.6 1 106

Reporting Limit for DF =1;	W	TOTAL	NA	μg/L
ND means not detected at or	S	TOTAL	5.0	mg/Kg
above the reporting limit				8 8

*water samples are reported in µg/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, filter samples in µg/filter.

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of $0.45 \mu m$ filtered and acidified sample.

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

Angela Rydelius, Lab Manager

DHS ELAP Certification 1644



Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled:	04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received:	04/15/11
	Client Contact: Peter Cusack	Date Extracted:	04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/16/11-04/18/11

Total Extractable Petroleum Hydrocarbons with Silica Gel Clean-Up*

Extraction method: SW35			methods: SW8015B	n Sinca Gei Clean-Op [*]	w	ork Order:	1104453
Lab ID	Client ID	Matrix	TPH-Diesel (C10-C23)	TPH-Motor Oil (C18-C36)	DF	% SS	Comments
1104453-001A	B-1-15	S	210	160	1	101	e3/e1,e7
1104453-002A	B-2-15	S	ND	ND	1	105	
1104453-003A	B-3-15	S	ND	ND	1	100	
1104453-004A	B-4-15	S	ND	ND	1	100	
1104453-005A	B-5-15	S	ND	ND	1	105	
1104453-006A	B-6-15	S	ND	ND	1	101	
1104453-007A	B-7-15	S	45	120	1	106	e7,e2
1104453-008A	B-8-15	S	ND	ND	1	102	
1104453-009A	B-9-15	S	ND	ND	1	105	
	g Limit for DF =1;	W NA NA		ug/L			
	s not detected at or	S	1.0	5.0		mg/l	Kg

ı	ND 414 414	**	1471	1177	ug/L
	ND means not detected at or	S	1.0	5.0	mg/Kg
L	above the reporting limit	5	1.0	3.0	1119 115
Ī		, · · · · · · · · · · · · · · · · · · ·	, 1: 1/ 1 1 1 /	1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	1' '1 1 ' 7 1 11

^{*} water samples are reported in μ g/L, wipe samples in μ g/wipe, soil/solid/sludge samples in mg/kg, product/oil/non-aqueous liquid samples in mg/L, and all DISTLC / STLC / SPLP / TCLP extracts are reported in μ g/L.

%SS = Percent Recovery of Surrogate Standard. DF = Dilution Factor

- +The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:
- e2) diesel range compounds are significant; no recognizable pattern
- e3) aged diesel is significant; and/or e1) unmodified or weakly modified diesel is significant
- e7) oil range compounds are significant



[#] cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract.

QC SUMMARY REPORT FOR SW8270C

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57651 WorkOrder 1104453

EPA Method SW8270C Extraction SW3550B Spiked Sample ID: 1104350-002A									02A			
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	1
7 mary to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Acenaphthene	ND<16	2	NR	NR	NR	74.8	74.8	0	30 - 130	30	30 - 130	30
4-Chloro-3-methylphenol	ND<16	4	NR	NR	NR	83.5	83.6	0.0778	30 - 130	30	30 - 130	30
2-Chlorophenol	ND<16	4	NR	NR	NR	102	102	0	30 - 130	30	30 - 130	30
1,4-Dichlorobenzene	ND<16	2	NR	NR	NR	93.5	93.6	0.0748	30 - 130	30	30 - 130	30
2,4-Dinitrotoluene	ND<16	2	NR	NR	NR	87	86.5	0.657	30 - 130	30	30 - 130	30
4-Nitrophenol	ND<80	4	NR	NR	NR	74.6	70.4	5.73	30 - 130	30	30 - 130	30
N-Nitrosodi-n-propylamine	ND<16	2	NR	NR	NR	120	120	0	30 - 130	30	30 - 130	30
Pentachlorophenol	ND<80	4	NR	NR	NR	30.6	30.6	0	30 - 130	30	30 - 130	30
Phenol	ND<16	4	NR	NR	NR	91.5	91.3	0.246	30 - 130	30	30 - 130	30
Pyrene	ND<16	2	NR	NR	NR	67.7	67.6	0.0443	30 - 130	30	30 - 130	30
1,2,4-Trichlorobenzene	ND<16	2	NR	NR	NR	89.6	87.1	2.77	30 - 130	30	30 - 130	30
%SS1:	73	200	#	#	N/A	85	83	3.22	30 - 130	30	30 - 130	30
%SS2:	#	200	58	50	14.6	87	86	1.54	30 - 130	30	30 - 130	30
%SS3:	88	200	91	88	3.32	98	94	4.44	30 - 130	30	30 - 130	30
%SS4:	72	200	74	75	1.55	83	82	1.21	30 - 130	30	30 - 130	30
%SS5:	#	200	#	111	N/A	92	90	1.36	30 - 130	30	30 - 130	30
%SS6:	68	200	69	69	0	69	69	0	30 - 130	30	30 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57651 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104453-001A	04/15/11 1:50 PM	04/15/11	04/19/11 4:34 AM	1104453-002A	04/15/11 1:55 PM	04/15/11	04/18/11 6:18 PM
1104453-003A	04/15/11 2:00 PM	04/15/11	04/18/11 7:37 PM	1104453-004A	04/15/11 2:05 PM	04/15/11	04/18/11 8:55 PM
1104453-005A	04/15/11 2:10 PM	04/15/11	04/18/11 10:12 PM	1104453-006A	04/15/11 2:15 PM	04/15/11	04/18/11 11:29 PM
1104453-007A	04/15/11 2:20 PM	04/15/11	04/19/11 12:46 AM	1104453-008A	04/15/11 2:25 PM	04/15/11	04/19/11 2:03 AM
1104453-009A	04/15/11 2:30 PM	04/15/11	04/19/11 3:19 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and / or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix, sample diluted due to high matrix or analyte content, or MS/MSD samples diluted due to high organic content.

#) surrogate diluted out of range; & = low or no recovery of surrogate or target analytes due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.

QA/QC Officer

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

QC SUMMARY REPORT FOR SW8021B/8015Bm

QC Matrix: Soil BatchID: 57681 WorkOrder 1104453 W.O. Sample Matrix: Soil

EPA Method SW8015Bm Extraction SW5030B Spiked S											: 1104409-0	13A
Analyte	Sample	mple Spiked MS MSD MS-N					LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex [£]	ND	0.60	110	116	5.22	113	104	8.78	70 - 130	20	70 - 130	20
MTBE	ND	0.10	108	122	12.7	110	110	0	70 - 130	20	70 - 130	20
Benzene	ND	0.10	84.3	92.2	8.85	89.2	83.6	6.43	70 - 130	20	70 - 130	20
Toluene	ND	0.10	82.9	89.5	7.68	87.9	82.4	6.47	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	84.9	90.4	6.28	89.4	84.4	5.71	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	85.1	90.3	5.96	89.3	84.6	5.34	70 - 130	20	70 - 130	20
%SS:	85	0.10	80	81	1.11	83	84	1.46	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57681 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104453-001A	04/15/11 1:50 PM	04/15/11	04/15/11 11:55 PM	1104453-002A	04/15/11 1:55 PM	04/15/11	04/16/11 12:24 AM
1104453-003A	04/15/11 2:00 PM	04/15/11	04/16/11 2:24 AM	1104453-004A	04/15/11 2:05 PM	04/15/11	04/19/11 1:33 AM
1104453-005A	04/15/11 2:10 PM	04/15/11	04/16/11 3:23 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

A QA/QC Officer

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

QC SUMMARY REPORT FOR SW8021B/8015Bm

QC Matrix: Soil BatchID: 57720 WorkOrder 1104453 W.O. Sample Matrix: Soil

EPA Method SW8015Bm Extraction SW5030B Spiked Sample ID: 1104453-009A												
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%))
	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex [£]	ND	0.60	99.5	103	3.37	101	99.6	0.934	70 - 130	20	70 - 130	20
MTBE	ND	0.10	109	106	3.39	106	107	0.515	70 - 130	20	70 - 130	20
Benzene	ND	0.10	105	104	0.374	105	104	1.09	70 - 130	20	70 - 130	20
Toluene	ND	0.10	92.1	92.1	0	92	90.8	1.26	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	94.1	94.6	0.490	93.6	92.4	1.25	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	106	107	0.969	106	104	1.99	70 - 130	20	70 - 130	20
%SS:	97	0.10	108	106	2.74	105	106	1.03	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57720 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104453-006A	04/15/11 2:15 PM	04/15/11	04/16/11 4:22 AM	1104453-007A	04/15/11 2:20 PM	04/15/11	04/19/11 2:05 AM
1104453-008A	04/15/11 2:25 PM	04/15/11	04/16/11 5:22 AM	1104453-009A	04/15/11 2:30 PM	04/15/11	04/16/11 5:51 AM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

A QA/QC Officer

QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104453

EPA Method SW6010B			Extraction SW3050B B					BatchID: 57629 Spike			ed Sample ID: 1104409-013A		
Analyte	Sample	Spiked	MS	MSD	MS-MSD	Spiked	LCS	LCSD	LCS-LCSD	Acceptance Criteria (%)			5)
	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Lead	5.7	50	100	103	2.44	10	94.6	95.1	0.580	75 - 125	25	75 - 125	25
%SS:	99	500	97	99	2.40	500	103	106	2.68	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57629 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104453-001A	04/15/11 1:50 PM	04/15/11	04/18/11 2:03 PM	1104453-002A	04/15/11 1:55 PM	04/15/11	04/18/11 2:05 PM
1104453-003A	04/15/11 2:00 PM	04/15/11	04/18/11 2:07 PM	1104453-004A	04/15/11 2:05 PM	04/15/11	04/18/11 2:09 PM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer

QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104453

EPA Method SW6010B		Extract	ion SW3	3050B		BatchID): 57721	Spiked Sample ID: 1104453-009A					
Analyte	Analyte Sample Spiked MS MSD MS-				MS-MSD	Spiked	LCS	LCSD	LCS-LCSD	Accepta		ance Criteria (%)	
, mary to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Lead	7.6	50	92.5	97.7	4.76	10	97.4	106	8.12	75 - 125	25	75 - 125	25
%SS:	106	500	103	106	2.30	500	106	103	2.97	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57721 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104453-005A	04/15/11 2:10 PM	04/15/11	04/18/11 2:11 PM	1104453-006A	04/15/11 2:15 PM	04/15/11	04/18/11 2:13 PM
1104453-007A	04/15/11 2:20 PM	04/15/11	04/18/11 2:15 PM	1104453-008A	04/15/11 2:25 PM	04/15/11	04/18/11 2:17 PM
1104453-009A	04/15/11 2:30 PM	04/15/11	04/18/11 1:12 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer

QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57680 WorkOrder 1104453

EPA Method SW8015B	Spiked Sample ID: 1104409-013A											
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
, many to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	ND	40	128	127	0.810	110	111	0.521	70 - 130	30	70 - 130	30
%SS:	101	25	107	109	2.16	94	96	1.39	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57680 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104453-001A	04/15/11 1:50 PM	04/15/11	04/16/11 8:13 AM	1104453-002A	04/15/11 1:55 PM	04/15/11	04/18/11 5:59 PM
1104453-003A	04/15/11 2:00 PM	04/15/11	04/16/11 5:53 AM	1104453-004A	04/15/11 2:05 PM	04/15/11	04/16/11 7:03 AM
1104453-005A	04/15/11 2:10 PM	04/15/11	04/16/11 10:41 AM	1104453-006A	04/15/11 2:15 PM	04/15/11	04/16/11 11:57 AM

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

DHS ELAP Certification 1644 QA/QC Officer

QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57719 WorkOrder 1104453

EPA Method SW8015B	Spiked Sample ID: 1104453-009A											
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%))
7 many to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	ND	40	119	121	2.06	93.7	94	0.340	70 - 130	30	70 - 130	30
%SS:	105	25	101	103	1.38	99	99	0	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57719 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104453-007A	04/15/11 2:20 PM	04/15/11	04/16/11 1:13 PM	1104453-008A	04/15/11 2:25 PM	04/15/11	04/16/11 4:58 PM
1104453-009A	04/15/11 2:30 PM	04/15/11	04/16/11 6:14 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

DHS ELAP Certification 1644

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer

McCampbell Analytical,	Inc.
"When Quality Counts"	

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled:
555 Montgomery St., Suite 1300		Date Received:
555 Mongomery Bu, Bulle 1500	Client Contact: Peter Cusack	Date Reported: 04/19/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/25/11

WorkOrder: 1104453

Enclosed within are:

- 1) The results of the analyzed samples from your project:
- 2) A QC report for the above samples,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.

1104453



Treadwell&Rollo

CHAIN OF CUSTODY RECORD

Page of

Environmental and Geote	chnical Consu	ltant																				5.95	55.9041		1	30	
		Firepo															500/F					16.5	565.7412				
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Job Number:	730 4	82302	0 (N.		Ar	naly	sis R	equ	este	d					Turna	round	
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Field Sample Identification No.	Date	Time	Lab Sample No.	Soil	Water	Other	HCL	H2SO2	HNO3	Other		5	H	60	1	METGY					Silica del	Diela Bala	Hold	R	emarks		
B-1-15	4-15-11	1350		X		\exists		1	λ			X	k	\vee	C	X			\Box		Y	_					
B-2-15	1	1355		7		\exists	\top	\top	X			X	X	X			\Box	\top	П	\top	Ü	,					
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B-4-15		1403		X		\exists		\top	X			X	X	2			\Box	\top		\top	1						
B-5-15		1410		8		\neg			X			7	Ċ	X			\Box	\top	1	\top	X						
B-6-15		1413	- X	N.					X		1	0	X	X			П		\Box		X						
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ICE/P GOOD CONDIT HEAD SPACE A DECHLORINAL	TON BSENT TED IN LAB VOAS O	White Copy APPROPRI CONTAINE PRESERVI	ERS		Yello	ow (Сору	y - L	.aboi	rator	y			8	P	ink (Сору -	Fiel	d		(CO	OC Number:	00	45 3	7	

1534 Willow Pass Rd Pittsburg C Pittsburg, CA 94565-1701

CHAIN-OF-CUSTODY RECORD

Page 1 of 1

Pittsburg (925) 25	g, CA 94565-1701 52-9262				V	VorkO	rder:	110445	53 A		Client(Code: 1	ſWRF				
		WaterTrax	Write	On EDF		Excel	[Fax	[✓ Email		Hard	Сору	Thi	rdParty	☐ J-1	flag
San Francisc		cc: PO:		eadwellrollo.com 5812 Hollis St.		l	Treadwell 555 Montg			nts Payable vell & Rollo intgomery St., Suite 130 ancisco, CA 94111			Requested TAT: Date Received: Date Add-On: Date Printed:			•	
									Req	uested	Tests	(See le	jend bo	elow)			-
Lab ID	Client ID		Matrix	Collection Date	Hold	1	2	3	4	5	6	7	8	9	10	11	12
1104453-001	B-1-15		Soil	4/15/2011 13:50		В											

Test Legend:

1 G-MBTEX_S	2	3	4	5
6	7	8	9	10
11	12]		

Prepared by: Maria Venegas

Comments: SEND HARD COPY. 72hr Rush. MBTEX added on 001 4/25/11 24hr

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled:	04/15/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received:	04/15/11
	Client Contact: Peter Cusack	Date Extracted:	04/15/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/15/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline with BTEX and MTBE*

Extraction	method: SW5030B		unge (ev erz)	-	tical methods:				Wor	k Order:	1104453
Lab ID	Client ID	Matrix	TPH(g)	MTBE	Benzene	Toluene	Ethylbenzene	Xylenes	DF	% SS	Comments
001A	B-1-15	S	5.2	ND	ND	ND	ND	ND	1	96	d7
	ing Limit for DF =1;	W	50	5.0	0.5	0.5	0.5	0.5		ug/I	
	ans not detected at or the reporting limit	S	1.0	0.05	0.005	0.005	0.005	0.005		mg/K	K g

* water and vapor samples are reported in µg/L, soil/sludge/solid samples in mg/kg,	wipe samples in µg/wipe, product/oil/non-aqueous liquid samples and all
TCLP & SPLP extracts in mg/L.	

- +The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:
- d7) strongly aged gasoline or diesel range compounds are significant in the TPH(g) chromatogram

____Angela Rydelius, Lab Manager

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57681 WorkOrder 1104453

EPA Method SW8021B/8015Bm	Extrac	ction SW	5030B					S	Spiked San	nple ID:	: 1104409-0	13A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
7 thaty to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex)	ND	0.60	110	116	5.22	113	104	8.78	70 - 130	20	70 - 130	20
MTBE	ND	0.10	108	122	12.7	110	110	0	70 - 130	20	70 - 130	20
Benzene	ND	0.10	84.3	92.2	8.85	89.2	83.6	6.43	70 - 130	20	70 - 130	20
Toluene	ND	0.10	82.9	89.5	7.68	87.9	82.4	6.47	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	84.9	90.4	6.28	89.4	84.4	5.71	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	85.1	90.3	5.96	89.3	84.6	5.34	70 - 130	20	70 - 130	20
%SS:	85	0.10	80	81	1.11	83	84	1.46	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57681 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104453-001A	04/15/11 1:50 PM	04/15/11	04/15/11 11:55 PM	1104453-002A	04/15/11 1:55 PM	04/15/11	04/16/11 12:24 AM
1104453-003A	04/15/11 2:00 PM	04/15/11	04/16/11 2:24 AM	1104453-004A	04/15/11 2:05 PM	04/15/11	04/19/11 1:33 AM
1104453-005A	04/15/11 2:10 PM	04/15/11	04/16/11 3:23 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

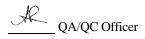
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



McCampbell Analytical, Inc.
"When Quality Counts"

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300		Date Received: 04/18/11
555 Mongomery Bu, Bulle 1500	Client Contact: Peter Cusack	Date Reported: 04/21/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/21/11

WorkOrder: 1104506

April 21, 2011

1100*	Peter:
Dear	reiei.

Enclosed within are:

- 1) The results of the 7 analyzed samples from your project: #730482302; 5812 Hollis St.,
- 2) A QC report for the above samples,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.



Treadwell&Rollo CHAIN OF CUSTODY RECORD

Number:	73040	2 - 4 - 1		_										Anal	ysis I	Requ	ested				Turnaround
ect Manager\Co plers:	ntact:	Peter	Cusack	_		_					1										72 Time
order (Signature	Required):	4	11)	- 3			No	. Co	ntai	ners	7	15							clean-up		400
				N	latri	_		_	serv	ative		2	H						cles		
ield Sample ntification No.	Date	Time	Lab Sample No.	Soil	Water	Other	HCL Hees	ONH ONH	loe	Other	HOT	SV	60						Silica gel	PloH	Remarks
10-15	4-18-11	(205		K					X		X	X	X						K)		
11-15		1210		X	1 1				X		X	1	X						X		
15-15		1218		X	*.				K		X	X	X					1			0 =
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15-15		1230		X	Н	+	+	+	X		1	X	X,	+	+		++		8		
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ished by: (Signatu	re)		Date () 18 11	<u> </u>			Time!	41			Red	eive	d by:	(Signat	ure)			D	ate	1 - 1 - 9	Time
ished by: (Signatu	re)		Date			7	Time				Rec			Lab: (Si		5	7	D	ate/	18/11	Time / 6/5
Laboratory (Natory Commen		McCo	impbell								Me	_		Shipm Carried	_		Lab co Courier		Vam	Fed Ex	Airborne l

PRESERVATION

1534 Willow Pass Rd Pittsburg, CA 94565-1701

CHAIN-OF-CUSTODY RECORD

ClientCode: TWRF

WorkOrder: 1104506

Page 1 of 1

(923) 232-9202	WaterTrax	WriteOn	☐ EDF	Excel	Fax	✓ Email	HardCopy	ThirdParty	J-flag
Report to:				Bi	II to:		Rec	uested TAT:	3 days
Peter Cusack Treadwell & Rollo 555 Montgomery St., Suite 1300	Email: p cc: PO:	ojcusack@treadv	vellrollo.com		Accounts Pa Treadwell & 555 Montgoi	,	00 Da i	te Received:	04/18/2011
San Francisco, CA 94111 (415) 955-5244 FAX (415) 955-9041		#730482302; 58 ²	12 Hollis St.		•	sco, CA 94111		te Printed:	04/18/2011

								R	Requ	ested	Tests (See le	gend be	elow)			
Lab ID	Client ID	Matrix	Collection Date	Hold	1	2	3	4	4	5	6	7	8	9	10	11	12
1104506-001	B-10-15	Soil	4/18/2011 12:05		Α	Α	Α										
1104506-002	B-11-15	Soil	4/18/2011 12:10		Α	Α	Α										
1104506-003	B-12-15	Soil	4/18/2011 12:15		Α	Α	Α										
1104506-004	B-13-15	Soil	4/18/2011 12:20		Α	Α	Α										
1104506-005	B-14-15	Soil	4/18/2011 12:25		Α	Α	Α										
1104506-006	B-15-15	Soil	4/18/2011 12:30		Α	Α	Α										
1104506-007	B-16-15	Soil	4/18/2011 12:00		Α	Α	Α										

Test Legend:				
1 8270D_S	2 PB_S	3 TPH(DMO)WSG_S	4	5
6	7	8	9	10
11	12			
The following SampIDs: 001A, 002A	, 003A, 004A, 005A, 006A, 007A contain t	estgroup.		Prepared by: Maria Venegas

Comments: SEND HARD COPY. 72hr Rush

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

Sample Receipt Checklist

Client Name:	Treadwell & R	(0110				Date a	and Time Received:	4/18/2011	4:15:41 PM
Project Name:	#730482302; 5	812 Hollis	St.			Check	dist completed and	reviewed by:	Maria Venegas
WorkOrder N°:	1104506	Matrix S	<u>soil</u>			Carrie	r: <u>Benjamin Ysl</u> a	as (MAI Courie	<u>r)</u>
			Chain	of Cu	stody (C	COC) Informa	ation		
Chain of custody	present?			Yes	V	No 🗆			
Chain of custody	signed when reli	nquished and r	eceived?	Yes	V	No 🗆			
Chain of custody	agrees with sam	ple labels?		Yes	✓	No 🗌			
Sample IDs noted	by Client on COC	?		Yes	V	No 🗆			
Date and Time of	collection noted b	y Client on CO	C?	Yes	✓	No 🗆			
Sampler's name no	oted on COC?			Yes	~	No 🗆			
			Sa	ample	Receipt	t Information	1		
Custody seals into	act on shipping co	ontainer/cooler		Yes		No 🗆		NA 🔽	
Shipping containe	er/cooler in good o	condition?		Yes	V	No 🗆			
Samples in prope	r containers/bottl	es?		Yes	~	No 🗆			
Sample containers	rs intact?			Yes	✓	No 🗆			
Sufficient sample	volume for indica	ited test?		Yes	✓	No 🗌			
		<u>Sam</u>	ple Preser	vatio	n and Ho	old Time (HT)) Information		
All samples receiv	ved within holding	time?		Yes	✓	No 🗌			
Container/Temp B	Blank temperature			Coole	er Temp:	5.8°C		NA 🗆	
Water - VOA vials	s have zero head	space / no but	obles?	Yes		No 🗆	No VOA vials subr	mitted 🗹	
Sample labels che	ecked for correct	preservation?		Yes	~	No 🗌			
Metal - pH accepta	able upon receipt	(pH<2)?		Yes		No 🗆		NA 🔽	
Samples Received	d on Ice?			Yes	✓	No 🗆			
			(Ice Type	e: WE	TICE)			
* NOTE: If the "No	lo" box is checke	d, see commei	nts below.						
	=====		====		===	====	=====		=====
Client contacted:		D	ate contact	ed:			Contacte	d by:	
Comments:									

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	,	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/18/11
	Client Contact: Peter Cusack	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104506

Lab ID	Lab ID 1104506-001A						
Client ID				B-10-15			
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surro	gate Re	coveries (%)			
<u> </u>					1		

Surrogate Recoveries (%)				
%SS1:	95	%SS2:	96	
%SS3:	91	%SS4:	81	
%SS5:	85	%SS6:	89	

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	•	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/18/11
	Client Contact: Peter Cusack	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Work Order: 1104506 Extraction Method: SW3550B

Lab ID		1104506-002A					
Client ID		B-11-15					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surro	ogate Re	coveries (%)			
%SS1:	82			%SS2:	89)	
% SS3·	87			%SS4·	82		

%SS3 %SS4

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/18/11
	Client Contact: Peter Cusack	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104506

Lab ID	1104506-003A						
Client ID		B-12-15					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surre	ogate Re	coveries (%)			
%SS1:	86			%SS2:	82		

Surrogate Recoveries (%)					
%SS1:	86	%SS2:	82		
%SS3:	92	%SS4:	78		
%SS5:	83	%SS6:	89		

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	3	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/18/11
	Client Contact: Peter Cusack	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104506

Lab ID	1104506-004A						
Client ID		B-13-15					
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
		Surre	ogate Re	coveries (%)			
%SS1:	89	-		%SS2:	88		

Surrogate Recoveries (%)					
%SS1:	89	%SS2:	88		
%SS3:	92	%SS4:	75		
%SS5:	69	%SS6:	87		
70000	· · · · · · · · · · · · · · · · · · ·	70.00	·		

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in $\mu g/L$, soil/sludge/solid samples in mg/kg, wipe samples in $\mu g/wipe$, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

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Treadwell & Rollo	•	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/18/11
	Client Contact: Peter Cusack	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104506

Lab ID	1104506-005A							
Client ID		B-14-15						
Matrix				Soil				
Compound	Concentration *	entration * DF Reporting Limit Compound Concentration *						
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
		Surre	gate Re	coveries (%)				
%SS1:	89			%SS2:	85	•		

Surrogate Recoveries (%)					
%SS1:	89	%SS2:	85		
%SS3:	84	%SS4:	77		
%SS5:	59	%SS6:	68		

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

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Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/18/11
	Client Contact: Peter Cusack	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Analytical Method: SW8270C Work Order: 1104506 Extraction Method: SW3550B

Lab ID	1104506-006A							
Client ID		B-15-15						
Matrix		Soil						
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33	, , , , , , , , , , , , , , , , , , , ,				
		Surro		coveries (%)				
%SS1:	94		g	%SS2:	81			
%SS3:	101			%SS4:	81			

Surrogate Recoveries (%)							
%SS1:	94	%SS2:	81				
%SS3:	101	%SS4:	83				
%SS5:	57	%SS6:	83				

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Lab ID

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/18/11
333 Wongomery St., State 1300	Client Contact: Peter Cusack	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/20/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

1104506-007A

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104506

Client ID		B-16-15							
				Soil					
Matrix			Ι	5011	T	1	L .		
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit		
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33		
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33		
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6		
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33		
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33		
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6		
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33		
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33		
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33		
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66		
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33		
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33		
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33		
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33		
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33		
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66		
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33		
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33		
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6		
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33		
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33		
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33		
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33		
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33		
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33		
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33		
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33		
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6		
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33		
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6		
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33		
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33		
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33		
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33		
2.4.6-Trichlorophenol	ND	1.0	0.33	, ,,,					
	1,2			coveries (%)					

Surrogate Recoveries (%)					
%SS1:	94	%SS2:	105		
%SS3:	96	%SS4:	83		
%SS5:	111	%SS6:	91		

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in $\mu g/L$, soil/sludge/solid samples in mg/kg, wipe samples in $\mu g/wipe$, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/18/11
	Client Contact: Peter Cusack	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 04/19/11-04/21/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline* Analytical methods SW8015Bm Extraction method SW5030B 1104506 Lab ID Client ID Matrix TPH(g) DF % SS Comments 001A B-10-15 S ND 002A S 102 B-11-15 ND 1 003A S 91 B-12-15 ND 1 004A B-13-15 S ND 1 89 B-14-15 S ď7 005A 3.5 1 91 006A B-15-15 S 21 1 89 d7 007A B-16-15 S ND 1 84

* water and vapor samples are reported in μg/L, soil/sludge/solid samples in mg/kg	, wipe samples in µg/wipe, product/oil/non-aqueous liquid
samples and all TCLP & SPLP extracts in mg/L.	

W

S

NA

1.0

d7) strongly aged gasoline or diesel range compounds are significant in the TPH(g) chromatogram

Angela Rydelius, Lab Manager

NA

mg/Kg

Reporting Limit for DF =1;

ND means not detected at or

above the reporting limit

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:



Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/18/11
	Client Contact: Peter Cusack	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/19/11

Lead by ICP*

Extraction method: SW30	50B	Analy	rtical methods: SW60	010B		Work Ord	der: 1104506
Lab ID	Client ID	Matrix	Extraction Type	Lead	DF	% SS	Comments
1104506-001A	B-10-15	S	TOTAL	ND	1	106	
1104506-002A	B-11-15	S	TOTAL	7.3	1	107	
1104506-003A	B-12-15	S	TOTAL	8.2	1	111	
1104506-004A	B-13-15	S	TOTAL	13	1	108	
1104506-005A	B-14-15	S	TOTAL	7.6	1	109	
1104506-006A	B-15-15	S	TOTAL	7.2	1	106	
1104506-007A	B-16-15	S	TOTAL	7.0	1	108	
						İ	†

Reporting Limit for DF =1;	W	TOTAL	NA	μg/L
ND means not detected at or	9	TOTAI	5.0	ma/Va
above the reporting limit	5	TOTAL	3.0	mg/Kg

*water samples are reported in µg/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, filter samples in µg/filter.

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of $0.45 \mu m$ filtered and acidified sample.

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

Angela Rydelius, Lab Manager

DHS ELAP Certification 1644



1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

	Client Project ID: #730482302; 5812	Date Sampled:	04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received:	04/18/11
	Client Contact: Peter Cusack	Date Extracted:	04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/19/11-04/21/11

Total Extractable Petroleum Hydrocarbons with Silica Gel Clean-Up*

Extraction method: SW3550B/3630C Analytical methods: SW8015B Work Order: 1104506 TPH-Diesel TPH-Motor Oil Client ID DF % SS Lab ID Matrix Comments (C10-C23) (C18-C36) 1104506-001A B-10-15 S ND ND 113 1104506-002A B-11-15 S ND ND 118 1104506-003A B-12-15 S 1.8 ND 117 1104506-004A B-13-15 S 10 ND 116 e1 1104506-005A B-14-15 S 52 44 111 1 e1 1104506-006A B-15-15 S 5 91 630 240 e1 1104506-007A B-16-15 S ND ND 112

Reporting Limit for DF =1;	W	NA	NA	ug/L
ND means not detected at or above the reporting limit	S	1.0	5.0	mg/Kg

^{*} water samples are reported in µg/L, wipe samples in µg/wipe, soil/solid/sludge samples in mg/kg, product/oil/non-aqueous liquid samples in mg/L, and all DISTLC / STLC / SPLP / TCLP extracts are reported in µg/L.

cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract.

%SS = Percent Recovery of Surrogate Standard. DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

e1) unmodified or weakly modified diesel is significant

e2) diesel range compounds are significant; no recognizable pattern

Angela Rydelius, Lab Manager

QC SUMMARY REPORT FOR SW8270C

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57743 WorkOrder 1104506

EPA Method SW8270C	Extra	ction SW	3550B					S	Spiked Sar	nple ID	: 1104506-0)07A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acc	eptance	Criteria (%)	
, undry to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Acenaphthene	ND	2	70.7	69	2.51	69.6	68.4	1.81	30 - 130	30	30 - 130	30
4-Chloro-3-methylphenol	ND	4	80.2	82.2	2.45	75.1	75.1	0	30 - 130	30	30 - 130	30
2-Chlorophenol	ND	4	70.8	67.9	4.17	75.4	70.2	7.10	30 - 130	30	30 - 130	30
1,4-Dichlorobenzene	ND	2	76.7	73.5	4.14	74.6	74.7	0.0804	30 - 130	30	30 - 130	30
2,4-Dinitrotoluene	ND	2	83.1	82.1	1.17	85	80.2	5.74	30 - 130	30	30 - 130	30
4-Nitrophenol	ND	4	72	71.3	1.05	66.3	69.8	5.17	30 - 130	30	30 - 130	30
N-Nitrosodi-n-propylamine	ND	2	71.8	69.2	3.68	91	79.5	13.5	30 - 130	30	30 - 130	30
Pentachlorophenol	ND	4	66.3	64.5	2.78	60.6	58	4.41	30 - 130	30	30 - 130	30
Phenol	ND	4	69.9	68.5	2.08	79.4	73.1	8.23	30 - 130	30	30 - 130	30
Pyrene	ND	2	88.6	82.2	7.45	87.2	83.6	4.19	30 - 130	30	30 - 130	30
1,2,4-Trichlorobenzene	ND	2	83	81.1	2.34	77.8	79.6	2.29	30 - 130	30	30 - 130	30
%SS1:	94	200	88	85	3.77	101	88	13.9	30 - 130	30	30 - 130	30
%SS2:	105	200	90	87	3.58	106	94	12.1	30 - 130	30	30 - 130	30
%SS3:	96	200	96	96	0	99	96	3.61	30 - 130	30	30 - 130	30
%SS4:	83	200	76	80	5.10	77	76	0.310	30 - 130	30	30 - 130	30
%SS5:	111	200	105	95	10.1	99	97	2.83	30 - 130	30	30 - 130	30
%SS6:	91	200	91	85	7.23	88	92	5.04	30 - 130	30	30 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57743 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104506-001A	04/18/11 12:05 PM	04/18/11	04/21/11 12:27 PM	1104506-002A	04/18/11 12:10 PM	04/18/11	04/21/11 1:45 PM
1104506-003A	04/18/11 12:15 PM	04/18/11	04/21/11 3:05 PM	1104506-004A	04/18/11 12:20 PM	04/18/11	04/21/11 3:43 PM
1104506-005A	04/18/11 12:25 PM	04/18/11	04/21/11 3:28 PM	1104506-006A	04/18/11 12:30 PM	04/18/11	04/21/11 4:25 PM
1104506-007A	04/18/11 12:00 PM	04/18/11	04/20/11 1:03 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and / or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix, sample diluted due to high matrix or analyte content, or MS/MSD samples diluted due to high organic content.

#) surrogate diluted out of range; & = low or no recovery of surrogate or target analytes due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.

QA/QC Officer

QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57720 WorkOrder 1104506

EPA Method SW8015Bm	Extrac	ction SW	5030B					S	Spiked San	nple ID	: 1104453-0	09A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
7 mary to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex [£]	ND	0.60	99.5	103	3.37	101	99.6	0.934	70 - 130	20	70 - 130	20
MTBE	ND	0.10	109	106	3.39	106	107	0.515	70 - 130	20	70 - 130	20
Benzene	ND	0.10	105	104	0.374	105	104	1.09	70 - 130	20	70 - 130	20
Toluene	ND	0.10	92.1	92.1	0	92	90.8	1.26	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	94.1	94.6	0.490	93.6	92.4	1.25	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	106	107	0.969	106	104	1.99	70 - 130	20	70 - 130	20
%SS:	97	0.10	108	106	2.74	105	106	1.03	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57720 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104506-001A	04/18/11 12:05 PM	I 04/18/11	04/19/11 3:30 AM	1104506-002A	04/18/11 12:10 PM	04/18/11	04/19/11 4:29 AM
1104506-003A	04/18/11 12:15 PM	04/18/11	04/19/11 4:59 AM	1104506-004A	04/18/11 12:20 PM	04/18/11	04/19/11 11:52 PM
1104506-005A	04/18/11 12:25 PM	04/18/11	04/21/11 2:33 AM	1104506-006A	04/18/11 12:30 PM	04/18/11	04/19/11 6:01 PM
1104506-007A	04/18/11 12:00 PM	I 04/18/11	04/20/11 5:21 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

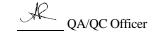
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104506

EPA Method SW6010B Extraction SW3050B					BatchID: 57721 Spiked Sample ID: 1104453-009A					9A			
Analyte	Sample	Spiked	MS	MSD	MS-MSD	Spiked	LCS	LCSD	LCS-LCSD	Acc	eptanc	e Criteria (%	5)
, inaly to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Lead	7.6	50	92.5	97.7	4.76	10	97.4	106	8.12	75 - 125	25	75 - 125	25
%SS:	106	500	103	106	2.30	500	106	103	2.97	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57721 SUMMARY

Lab ID	Date Sampled	Date Extracte	ed Date Analyzed	Lab ID	Date Sampled	Date Extracte	d Date Analyzed
1104506-001A	04/18/11 12:05 PM	04/18/11	04/19/11 11:35 AM	1104506-002A	04/18/11 12:10 PM	04/18/11	04/19/11 11:37 AM
1104506-003A	04/18/11 12:15 PM	04/18/11	04/19/11 11:39 AM	1104506-004A	04/18/11 12:20 PM	04/18/11	04/19/11 11:41 AM
1104506-005A	04/18/11 12:25 PM	04/18/11	04/19/11 11:43 AM	1104506-006A	04/18/11 12:30 PM	04/18/11	04/19/11 11:45 AM
1104506-007A	04/18/11 12:00 PM	04/18/11	04/19/11 11:47 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer

QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57719 WorkOrder 1104506

EPA Method SW8015B	W8015B Extraction SW3550B/3630C									nple ID:	: 1104453-0	009A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acc	eptance	Criteria (%)	١
, many to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	ND	40	119	121	2.06	93.7	94	0.340	70 - 130	30	70 - 130	30
%SS:	105	25	101	103	1.38	99	99	0	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57719 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104506-001A	04/18/11 12:05 PM	04/18/11	04/21/11 3:25 PM	1104506-002A	04/18/11 12:10 PM	04/18/11	04/19/11 11:16 AM
1104506-003A	04/18/11 12:15 PM	04/18/11	04/21/11 2:36 PM	1104506-004A	04/18/11 12:20 PM	04/18/11	04/19/11 10:06 AM
1104506-005A	04/18/11 12:25 PM	04/18/11	04/19/11 2:55 AM	1104506-006A	04/18/11 12:30 PM	04/18/11	04/19/11 4:07 PM
1104506-007A	04/18/11 12:00 PM	04/18/11	04/20/11 4:23 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer

DHS ELAP Certification 1644

McCampbell Analytical,	Inc.
"When Quality Counts"	

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled:
555 Montgomery St., Suite 1300		Date Received:
tee mongomery on, same 1999	Client Contact: Peter Cusack	Date Reported: 04/21/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/25/11

WorkOrder: 1104506

Enclosed within are:

- 1) The results of the analyzed samples from your project:
- 2) A QC report for the above samples,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.

1104506



Treadwell&Rollo Environmental and Geotechnical Consultant

CHAIN OF CUSTODY RECORD

555 Montgomery Street, Suite 1300, San Francisco, CA 94111 Ph: 415.955.9040/Fax: 415.955.9041

Page of

Site Name:	5812	Holls								cland (e 200,												6.56	5.7412				
Job Number:	7304	82307	7												Ana	lys	is Re	que	sted		H		1		Turnaro	ind	
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		- 04		N	latri	K				ative	1	1	7	2	7	,					clean-up		[_
Field Sample Identification No.	Date	Time	Lab Sample No.	Soil		_	H-SO	_		Other	TOH-		0/1-	189	MRTEX 4/26						Silica gel			Ren	narks		
B-10-15	4-18-11	1205		X					X		X	1	A Y								X						
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B-15-12		1218		K	*.	1	_		K		X	X	X								X						
13-13-15		1220		0	4	+	+	-	X		X	1				Ш	4	\perp		_	X						
B-14-15		1225		X	+	+	+	\vdash	X		X		X	9	+	\Box	+	-	1	+	V						
B-16-15*	M	1200		X	+	+	+	+	X		1x	44		4	+	H		+	-	+	1						-
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* Tube La	beled	White Copy	- Original	,	Yello	w C	ору -	- La	bora	atory]						Py - I			A			Number:	005	242		

DECILORINATED IN LAB

TRESERVATION

VOAS | O&G | METALS | OTHER

1534 Willow Pass Rd Pittsburg, CA 94565-1701

CHAIN-OF-CUSTODY RECORD

Page 1 of 1

(925) 252	-9262				`	VorkC	rder:	110450)6 A	Clien	tCode:	ľWRF				
		WaterTrax	Write	On EDF		Exce	[Fax	VE	imail	Hard	Сору	∏Thi	rdParty	☐ J-f	lag
	k Rollo cc: omery St., Suite 1300 PO: sco, CA 94111 ProjectNo: #730482302; 5812 Hollis St.						Tre 55:	eadwell 5 Montg	Payable & Rollo gomery St., sisco, CA 9		00	Dai Dai	juested te Rec te Add te Prin	eived: -On:	3 04/18/ 04/25/ 04/25/	/2011
									Reques	ted Tests	(See le	gend b	elow)			
Lab ID	Client ID		Matrix	Collection Date	Hold	1	2	3	4	5 6	7	8	9	10	11	12
1104506-002	B-11-15		Soil	4/18/2011 12:10		В										

Test Legend:

1 G-MBTEX_S	2	3	4	5
6	7	8	9	10
11	12			

Prepared by: Maria Venegas

Comments: SEND HARD COPY. 72hr Rush. MBTEX added on 002 4/25/11 24hr

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled:	04/18/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received:	04/18/11
	Client Contact: Peter Cusack	Date Extracted:	04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/19/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline with BTEX and MTBE*

	G	asomic 1	Mange (Co-C12)	voiame my	ui ocai bolis	as Gasoniic	WILLDIEAC	mu wiidi					
Extraction method: SW5030B Analytical methods: SW8021B/8015Em Work Order: 1104506 Lab ID Client ID Matrix TPH(g) MTBE Benzene Toluene Ethylbenzene Xylenes DF % SS Comments													
Lab ID	ab ID Client ID Matrix TPH(g) MTBE Benzene Toluene Ethylbenzene Xylenes DF % SS Comme												
002A	B-11-15	S	ND	ND	ND	ND	ND	ND	1	102			
	ng Limit for DF =1;	W	50	5.0	0.5	0.5	0.5	0.5		ug/L			
	ns not detected at or the reporting limit	S	1.0	0.05	0.005	0.005	0.005	0.005		mg/K	<u>Ig</u>		

^{*} water and vapor samples are reported in µg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts in mg/L.

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57720 WorkOrder 1104506

EPA Method SW8021B/8015Bm	Extrac	tion SW	5030B					S	Spiked San	nple ID	: 1104453-0	09A			
Analyte	Sample	Spiked MS MSD MS-MSD LCS LCSD LCS-LCSD Acceptance Criteria (%)													
7 tildiyto	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD			
TPH(btexf)	ND	0.60	99.5	103	3.37	101	99.6	0.934	70 - 130	20	70 - 130	20			
MTBE	ND	0.10	109	106	3.39	106	107	0.515	70 - 130	20	70 - 130	20			
Benzene	ND	0.10	105	104	0.374	105	104	1.09	70 - 130	20	70 - 130	20			
Toluene	ND	0.10	92.1	92.1	0	92	90.8	1.26	70 - 130	20	70 - 130	20			
Ethylbenzene	ND	0.10	94.1	94.6	0.490	93.6	92.4	1.25	70 - 130	20	70 - 130	20			
Xylenes	ND	0.30	106	107	0.969	106	104	1.99	70 - 130	20	70 - 130	20			
%SS:	97	0.10	108	106	2.74	105	106	1.03	70 - 130	20	70 - 130	20			

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57720 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104506-001A	04/18/11 12:05 PM	I 04/18/11	04/19/11 3:30 AM	1104506-002A	04/18/11 12:10 PM	04/18/11	04/19/11 4:29 AM
1104506-003A	04/18/11 12:15 PM	04/18/11	04/19/11 4:59 AM	1104506-004A	04/18/11 12:20 PM	04/18/11	04/19/11 11:52 PM
1104506-005A	04/18/11 12:25 PM	04/18/11	04/21/11 2:33 AM	1104506-006A	04/18/11 12:30 PM	04/18/11	04/19/11 6:01 PM
1104506-007A	04/18/11 12:00 PM	I 04/18/11	04/20/11 5:21 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

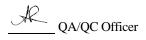
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



McCampbell Analytical,	Inc.
"When Ovelity Counts"	

Treadwell & Rollo	Client Project ID: #7304823021; 5812 Hollis St	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300		Date Received: 04/18/11
tee mongomery on, same 1999	Client Contact: Rob Milano	Date Reported: 04/25/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/22/11

WorkOrder: 1104507

April 25, 2011

\mathbf{r}		-	-		
11	ear	v.	\sim	h	
	ca	1	. , ,	,	

Enclosed within are:

- 1) The results of the 7 analyzed samples from your project: #7304823021; 5812 Hollis St,
- 2) A QC report for the above samples,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.

1104507

Treadwell&Rollo

Environmental and Geotechnical Consultant

CHAIN OF CUSTODY RECORD

555 Montgomery Street, Suite 1300, San Francisco, CA 94111 Ph: 415.955.9040/Fax: 415.955.9041

Page ___ of ___

	20.00		501 14 777 Ca						land C											16.5	565.	7412					
Site Name: Job Number: Project Manager\Cor Samplers: Recorder (Signature	Rob	Hollis 12302 Retes Miluna	Cusacic			N	o. Co	ontai	ners	-di-me	. 5	3	Ana						air dags	T				1	Turna	ne	
Field Sample Identification No.	Date	Time	Lab Sample No.	-	Water Other	-	H ₂ SO ₄	_	Other	TPH-a.	5002	100							lo loo coilio	D C	Hold			Re	marks		
P10-B P10-9 P10-10 P10-12 P10-13 P10-14	4-18-11 4-18-11 4-18-11 4-15-11 4-8-11 4-8-11	1235 1240 1245 1335 1330 1650 1645		X X X X X X				XXXXXX		XXXXXX	XXXXXX	XXXXXX							XXXXXX								
Relinquished by: (Signation of Signation	ure)	McC	Date 4-18-14 Date 18-14	(_	Time	618	1503		Red	ceive	ed by	(Sign	Signa) ature	0		Scour	Da	ate ate	4/1	18/1	//	Tir	ne /6	15	JPS
Sent to Laboratory (Laboratory Commer		McCa White Copy		,	Yellov	v Cop	oy - L	.abor	atory	Me		ICE/ GOO HEA DEC	° V	nk O	DPY ON BSEN	- Fie	ld B	Al	PPRO	PRI INE RVE	DC	N LAB			524		JPS

1534 Willow Pass Rd Pittsburg, CA 94565-1701

CHAIN-OF-CUSTODY RECORD

Page 1 of 1

(925) 252-9262				WorkOr	der: 110450'	7 Client	Code: TWRF		
	WaterTrax	WriteOn	☐ EDF	Excel	Fax	✓ Email	HardCopy	ThirdParty	J-flag
Report to:				Bil	I to:		Req	uested TAT:	5 days
Rob Milano	Email: rı	nmilano@tread\	wellrollo.com		Accounts Pa	yable			
Treadwell & Rollo	cc:				Treadwell &	Rollo	_		
555 Montgomery St., Suite 1300	PO:				555 Montgor	nery St., Suite 130	₀₀ Dai	e Received:	04/18/2011
San Francisco, CA 94111	ProjectNo: #	7304823021; 58	12 Hollis St		San Franciso	co, CA 94111	Dat	e Printed:	04/18/2011
(415) 955-5244 FAX (415) 955-9041									
						Requested Tests	(See legend b	elow)	

					Requested Tests (See legend below)											
Lab ID	Client ID	Matrix	Collection Date	Hold	1	2	3	4	5	6	7	8	9	10	11	12
1104507-001	P10-8	Soil	4/18/2011 12:35	ПП	A	Α	Α									
1104507-002	P10-9	Soil	4/18/2011 12:40		Α	Α	Α									
1104507-003	P10-10	Soil	4/18/2011 12:45		Α	Α	Α									
1104507-004	P10-11	Soil	4/18/2011 13:35		Α	Α	Α									
1104507-005	P10-12	Soil	4/18/2011 13:30		Α	Α	Α									
1104507-006	P10-13	Soil	4/18/2011 16:50		Α	Α	Α									
1104507-007	P10-14	Soil	4/18/2011 16:45		Α	Α	Α									

<u>Test Legend</u> :				
1 8270D_S	2 PB_S	3 TPH(DMO)WSG_S	4	5
6	7	8	9	10
11	12			
The following SampIDs: 001A, 002A	Prepared by: Ana Venegas			

Comments: SEND HARD COPY

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

Sample Receipt Checklist

Client Name:	Treadwell & Ro	ollo				Date	e ar	nd Time Received:	4/18/2011	4:56:11 PM
Project Name:	#730482302I; 5	812 Hollis	St			Che	ckl	list completed and re	viewed by:	Ana Venegas
WorkOrder N°:	1104507	Matrix S	<u>soil</u>			Carr	rier	r: <u>Benjamin Yslas</u>	(MAI Courie	<u>er)</u>
			Chain	of Cu	stody (C	COC) Inforn	nat	<u>tion</u>		
Chain of custody	y present?			Yes	V	No 🗆]			
Chain of custody	y signed when relin	quished and ı	eceived?	Yes	V	No 🗆]			
Chain of custody	y agrees with samp	le labels?		Yes	✓	No 🗆]			
Sample IDs noted	d by Client on COC?			Yes	V	No 🗆]			
Date and Time of	f collection noted by	Client on CO	C?	Yes	✓	No 🗆]			
Sampler's name	noted on COC?			Yes	✓	No 🗆]			
			Sa	ample	Receipt	t Informatio	<u>on</u>			
Custody seals in	itact on shipping co	ntainer/coole	?	Yes		No 🗆]		NA 🔽	
Shipping contain	ner/cooler in good co	ondition?		Yes	V	No 🗆]			
Samples in prop	er containers/bottle	s?		Yes	V	No 🗆]			
Sample containe	ers intact?			Yes	✓	No 🗆]			
Sufficient sample	e volume for indicat	ed test?		Yes	✓	No 🗌]			
		Sam	ple Preser	rvatio	n and Ho	old Time (H	IT)	Information		
All samples rece	eived within holding	time?		Yes	V	No 🗌]			
Container/Temp	Blank temperature			Coole	er Temp:	5.8°C			NA 🗆	
Water - VOA via	ıls have zero heads	pace / no bul	obles?	Yes		No 🗆]	No VOA vials submit	ted 🗸	
Sample labels cl	hecked for correct p	reservation?		Yes	✓	No 🗌]			
Metal - pH accep	otable upon receipt	(pH<2)?		Yes		No 🗆]	I	NA 🔽	
Samples Receive	ed on Ice?			Yes	✓	No \square]			
			(Ice Type	e: WE	TICE)				
* NOTE: If the "I	No" box is checked	, see comme	nts below.							
		====								
Client contacted:	:	D	ate contact	ed:				Contacted I	by:	
Comments:										

Treadwell & Rollo	Client Project ID: #7304823021; 5812	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/18/11
	Client Contact: Rob Milano	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104507

Lab ID	1104507-001A							
Client ID	P10-8							
Matrix		Soil						
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
		Surro	gate Re	coveries (%)	-			
%SS1:	85			%SS2:	82			
	-							

Surrogate Recoveries (%) %SS1: 85 %SS2: 82 %SS3: 84 %SS4: 77 %SS5: 70 %SS6: 78

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

McCampbell Analytical, Inc. "When Quality Counts"

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302l; 5812	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/18/11
	Client Contact: Rob Milano	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104507

Lab ID	1104507-002A							
Client ID		P10-9						
Matrix		Soil						
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
		Surre	gate Re	coveries (%)				
%SS1:	89			%SS2:	84			

Surrogate Recoveries (%)						
%SS1:	89	%SS2:	84			
%SS3:	87	%SS4:	78			
%SS5:	65	%SS6:	76			

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302l; 5812	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/18/11
	Client Contact: Rob Milano	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104507

Acetochlor	Lab ID		1104507-003A						
Compound Concentration DF Reporting Compound Concentration DF Reporting Lamba	Client ID	P10-10							
Compound Concentration	Matrix		Soil						
Acetochlor	Compound	Concentration *	DF		Compound	Concentration *	DF		
Benzidine	Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Benzo(a)anthracene ND 1.0 0.33 Benzo(b)fluoranthene ND 1.0 0.33 Benzo(a)fluoranthene ND 1.0 0.33 Benzo(g,h,i)perylene ND 1.0 0.33 Benzo(a)pyrene ND 1.0 0.33 Benzo(g,h,i)perylene ND 1.0 0.33 Benzo(a)pyrene ND 1.0 0.33 Benzo(g,h,i)perylene ND 1.0 0.33 Bis (2-chloroethy) Ether ND 1.0 0.33 Bis (2-chloroethoxy) Methane ND 1.0 0.33 Bis (2-chloroethyl) Ether ND 1.0 0.33 Bis (2-chloroethoxy) Methane ND 1.0 0.33 Bis (2-chlyhlexyl) Phthalate ND 1.0 0.33 4-Chloroanthene ND 1.0 0.33 Bis (2-chlyhlexyl) Phthalate ND 1.0 0.33 4-Chloroanthene ND 1.0 0.36 4-Chloro-3-methylphenol ND 1.0 0.33 4-Chloroanthene ND 1.0 0.36 4-Chloro-3-methylphenol ND 1.0 0.33 2-Chloroanthene ND 1.0 0.33 Chrysene ND 1.0 0.33 Dibenzo(a,h)anthracene ND 1.0 0.33 Dibenzofuran ND 1.0 0.33 Dibenzo(a,h)anthracene ND 1.0 0.33 Dibenzofuran ND 1.0 0.33 Di-n-butyl Phthalate ND 1.0 0.33 1,4-Dichlorobenzene ND 1.0 0.33 3,3-Dichlorobenzene ND 1.0 0.35 1,4-Dichlorophenol ND 1.0 0.33 3,3-Dichlorobenzene ND 1.0 0.35 1,4-Dimethylphenol ND 1.0 0.33 3,3-Dichlorobenzene ND 1.0 0.36 2,4-Dimethylphenol ND 1.0 0.33 3,3-Dichlorobenzene ND 1.0 0.35 1,4-Dimethylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.35 1,4-Dimethylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.35 1,4-Dimitro-2-methylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.	Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzo(k)fluoranthene ND	Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)pyrene ND	Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
1,1-Bipheny ND	Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Bis (2-chloroethyl) Ether ND 1.0 0.33 Bis (2-chloroisopropyl) Ether ND 1.0 0.33 Bis (2-chlyhexyl) Phthalate ND 1.0 0.33 4-Bromophenyl Phenyl Ether ND 1.0 0.33 Bis (2-chlyhexyl) Phthalate ND 1.0 0.33 4-Chloroaniline ND 1.0 0.66 4-Chloro-3-methylphenol ND 1.0 0.33 4-Chloroaniline ND 1.0 0.33 2-Chlorophenyl Phenyl Ether ND 1.0 0.33 1.0 0.34 0.35 0.	Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
Bis (2-ethylhexyl) Phthalate	1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Butylbenzyl Phthalate	Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
4-Chloro-3-methylphenol ND 1.0 0.33 2-Chloronaphthalene ND 1.0 0.33 2-Chlorophenol ND 1.0 0.33 4-Chlorophenyl Phenyl Ether ND 1.0 0.33 Chrysene ND 1.0 0.33 Dibenzo(a,h)anthracene ND 1.0 0.33 Dibenzo(aran ND 1.0 0.33 Dibenzo(a,h)anthracene ND 1.0 0.33 1,2-Dichlorobenzene ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.33 1,3-Dichlorobenzene ND 1.0 0.33 1,4-Dichlorobenzene ND 1.0 0.33 1,3-Dichlorobenzidine ND 1.0 0.66 0.4-Dichlorophenol ND 1.0 0.33 Diethyl Phthalate ND 1.0 0.63 0.33 0.4-Dichlorobenzene ND 1.0 0.33 Diethyl Phthalate ND 1.0 0.33 0.33 0.34-Dichlorobenzene ND 1.0 0.34-Dichlorobenzene ND 1.0 0.35 0.34-Dichlorobenzene ND 1.0 0.35 0.34-Dichlorobenzene ND 1.0 0.35 0.34-Dichlorobenzene ND 1.0 0.35 0.34-Dichloro	Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
2-Chlorophenol ND	Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
Chrysene	4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
Dibenzofuran ND	2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
1,2-Dichlorobenzene	Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
1,4-Dichlorobenzene ND 1.0 0.33 3,3-Dichlorobenzidine ND 1.0 0.66 2,4-Dichlorophenol ND 1.0 0.33 Diethyl Phthalate ND 1.0 0.33 2,4-Dimethylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 4,6-Dinitro-2-methylphenol ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Fluoranthene ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0	Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
2,4-Dichlorophenol ND 1.0 0.33 Diethyl Phthalate ND 1.0 0.33 2,4-Dimethylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 4,6-Dinitro-2-methylphenol ND 1.0 1.6 2,4-Dinitrophenol ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 1.6 Hexachlorocyclopentadiene ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylnaphthalene ND 1.0	1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
2,4-Dimethylphenol ND 1.0 0.33 Dimethyl Phthalate ND 1.0 0.33 4,6-Dinitro-2-methylphenol ND 1.0 1.6 2,4-Dinitrophenol ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 1.6 Hexachlorocyclopentadiene ND 1.0 0.33 Indenot (1,2,3-cd) pyrene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 2-Methylphenol (o-Cresol) ND 1.0 0.33 2-Nitroaniline ND 1.0<	1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
4,6-Dinitro-2-methylphenol ND 1.0 1.6 2,4-Dinitrophenol ND 1.0 1.6 2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 1.6 Hexachlorobutadiene ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 1.6 Hexachlorocyclopentadiene ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 <	2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dinitrotoluene ND 1.0 0.33 2,6-Dinitrotoluene ND 1.0 0.33 Di-n-octyl Phthalate ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 1.6 Hexachlorobutadiene ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 1.6 Hexachloroethane ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylphenol (m,p-Cres ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6	2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
Di-n-octyl Phthalate ND 1.0 0.33 1,2-Diphenylhydrazine ND 1.0 0.33 Fluoranthene ND 1.0 0.33 Fluorene ND 1.0 0.33 Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 1.6 Hexachloroethane ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 2-Methylphenol (o-Cresol) ND 1.0 0.33 3 &/or 4-Methylphenol (m,p-Cres ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitroaniline ND 1.0 1.6	4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
Fluoranthene	2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Hexachlorobenzene ND 1.0 0.33 Hexachlorobutadiene ND 1.0 0.33 Hexachlorocyclopentadiene ND 1.0 0.33 Indeno (1,2,3-cd) pyrene ND 1.0 0.33 Isophorone ND 1.0 0.33 2-Methylnaphthalene ND 1.0 0.33 2-Methylphenol (o-Cresol) ND 1.0 0.33 3 &/or 4-Methylphenol (m,p-Cres ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1.2,4-Trichlorobenzene ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33 2.4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33 ND 1.0 0.33 0.33 0.33 1.2,4-Trichlorophenol ND 1.0 0.33 1.3 1.3 1.3 1.3 1.3 1.4 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.5 1.	Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Hexachlorocyclopentadiene	Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene ND	Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
2-Methylnaphthalene ND 1.0 0.33 2-Methylphenol (o-Cresol) ND 1.0 0.33 3 &/or 4-Methylphenol (m,p-Cres) ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1.2,4-Trichlorobenzene ND 1.0 0.33 2.4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33 </td <td>Hexachlorocyclopentadiene</td> <td>ND</td> <td>1.0</td> <td>1.6</td> <td>Hexachloroethane</td> <td>ND</td> <td>1.0</td> <td>0.33</td>	Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres ND 1.0 0.33 Naphthalene ND 1.0 0.33 2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Nitroaniline ND 1.0 1.6 3-Nitroaniline ND 1.0 1.6 4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
4-Nitroaniline ND 1.0 1.6 Nitrobenzene ND 1.0 0.33 2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitrophenol ND 1.0 1.6 4-Nitrophenol ND 1.0 1.6 N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
N-Nitrosodiphenylamine ND 1.0 0.33 N-Nitrosodi-n-propylamine ND 1.0 0.33 Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33 ND 1.0 0.33	4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
Pentachlorophenol ND 1.0 1.6 Phenanthrene ND 1.0 0.33 Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2,4,6-Trichlorophenol ND 1.0 0.33	2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
Phenol ND 1.0 0.33 Pyrene ND 1.0 0.33 1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33	N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
1,2,4-Trichlorobenzene ND 1.0 0.33 2,4,5-Trichlorophenol ND 1.0 0.33 2.4.6-Trichlorophenol ND 1.0 0.33	Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
2.4.6-Trichlorophenol ND 1.0 0.33	Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
	1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
Surrogate Recoveries (%)	2.4.6-Trichlorophenol	ND	1.0	0.33					
			Surr	ogate Re	coveries (%)				

Surrogate Recoveries (%)						
%SS1:	88	%SS2:	82			
%SS3:	85	%SS4:	78			
%SS5:	63	%SS6:	79			
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Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

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Treadwell & Rollo	9	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/18/11
	Client Contact: Rob Milano	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104507

Lab ID	1104507-004A						
Client ID		P10-11					
Matrix		Soil					
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND 1.0 0.33						
		Surro	gate Re	coveries (%)			
%SS1:	89			%SS2:	81		
%SS3:	84			%SS4:	77		
,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	0-7			/000			

Comments

%SS6:

66

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in $\mu g/L$, soil/sludge/solid samples in mg/kg, wipe samples in $\mu g/wipe$, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302l; 5812	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/18/11
	Client Contact: Rob Milano	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/22/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104507

Lab ID	1104507-005A							
Client ID		P10-12						
Matrix				Soil				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33	
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33	
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6	
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33	
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33	
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6	
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33	
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33	
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33	
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66	
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33	
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33	
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33	
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33	
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33	
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66	
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33	
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33	
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6	
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33	
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33	
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33	
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33	
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33	
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33	
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33	
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6	
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33	
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6	
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33	
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33	
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33	
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33	
2.4.6-Trichlorophenol	ND	1.0	0.33					
		Surre	gate Re	coveries (%)				
%SS1:	92			%SS2:	73	73		

Surrogate Recoveries (%) %SS1: 92 %SS2: 73 %SS3: 88 %SS4: 76 %SS5: 61 %SS6: 81

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	3	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/18/11
555 Workgomery St., State 1500	Client Contact: Rob Milano	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/22/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104507

Extraction Method: SW 3550B		Anai	yticai Meti	nod: SW8270C	work Ord	er: 110	14507
Lab ID		1104507-006A					
Client ID				P10-13			
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
	-	Surr	ogate Re	coveries (%)	-	· ·	
%SS1:	87			%SS2:	82		
%SS3:	85			%SS4:	77		

Surrogate Recoveries (%)					
%SS1:	87	%SS2:	82		
%SS3:	85	%SS4:	77		
%SS5:	64	%SS6:	78		

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302l; 5812	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/18/11
	Client Contact: Rob Milano	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104507

Lab ID 1104507-007A							
Client ID P10-14							
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33			-	
	-	Surre	ogate Re	coveries (%)			
%SS1:	90			%SS2:	78		

%SS1: 90 %SS2: 78 %SS3: 82 %SS4: 75 %SS5: 62 %SS6: 70

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Work Order: 1104507

Treadwell & Rollo	Client Project ID: #730482302l; 5812 Hollis St	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/18/11
	Client Contact: Rob Milano	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 04/20/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline* Extraction method SW5030B Analytical methods SW8015Bm

Extraction inclined 5 w.	70300	Allaryticar	inctious 5 w 6015Diii	****	ik Oluci.	1107307
Lab ID	Client ID	Matrix	TPH(g)	DF	% SS	Comments
001A	P10-8	S	ND	1	86	
002A	P10-9	S	6.2	1	94	d7
003A	P10-10	S	ND	1	82	
004A	P10-11	S	ND	1	85	
005A	P10-12	S	14	1	77	d7
006A	P10-13	S	ND	1	83	
007A	P10-14	S	ND	1	83	
	ing Limit for DF =1;	W	NA		NA	•
ND mea	ans not detected at or	S	1.0		mg/Kg	

ND means not detected at or above the reporting limit	S	1.0	mg/Kg
* water and vapor samples are reported in ug/L soil/sluc	lge/solid sample	es in mg/kg_wine samples in ug/wine_product	/oil/non-aqueous liquid

^{*} water and vapor samples are reported in µg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts in mg/L.

d7) strongly aged gasoline or diesel range compounds are significant in the TPH(g) chromatogram

Angela Rydelius, Lab Manager

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:



Extraction method: SW3050B

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #7304823021; 5812	Date Sampled: 04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/18/11
<i>y</i> ,	Client Contact: Rob Milano	Date Extracted: 04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/19/11

Lead by ICP*

Analytical methods: SW6010B

Lab ID Client ID Matrix Extraction Type Lead DF % SS Comments 1104507-001A S TOTAL 1 P10-8 11 108 1104507-002A P10-9 S TOTAL 99 1 109 1104507-003A S TOTAL P10-10 12 1 109 1104507-004A P10-11 S TOTAL 13 1 113 1104507-005A P10-12 S TOTAL 5.1 108 1104507-006A P10-13 S TOTAL 8.7 108 1104507-007A P10-14 S TOTAL 9.3 1 113

Reporting Limit for DF =1;	W	TOTAL	NA	μg/L
ND means not detected at or above the reporting limit	S	TOTAL	5.0	mg/Kg

*water samples are reported in μ g/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, filter samples in μ g/filter.

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of $0.45 \mu m$ filtered and acidified sample.

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

Angela Rydelius, Lab Manager

DHS ELAP Certification 1644

Work Order: 1104507



	Client Project ID: #730482302l; 5812	Date Sampled:	04/18/11
555 Montgomery St., Suite 1300	Hollis St	Date Received:	04/18/11
	Client Contact: Rob Milano	Date Extracted:	04/18/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/19/11-04/22/11

Total Extractable Petroleum Hydrocarbons with Silica Gel Clean-Up*

Extraction method: SW	3550B/3630C		W	ork Order:	1104507		
Lab ID	Client ID	Matrix	TPH-Diesel (C10-C23)	TPH-Motor Oil (C18-C36)	DF	% SS	Comments
1104507-001A	P10-8	S	1.8	ND	1	110	e2
1104507-002A	P10-9	S	100	52	1	101	e1
1104507-003A	P10-10	S	2.5	ND	1	102	e1
1104507-004A	P10-11	S	ND	ND	1	108	
1104507-005A	P10-12	S	28	13	1	92	e1
1104507-006A	P10-13	S	ND	ND	1	107	
1104507-007A	P10-14	S	ND	ND	1	104	
Reporting Limit for DF =1;		W	NA	NA		ug/	 L
ND mea	ns not detected at or	 		<u> </u>	+		

Reporting Limit for DF =1;	W	NA	NA	ug/L
ND means not detected at or above the reporting limit	S	1.0	5.0	mg/Kg

^{*} water samples are reported in μ g/L, wipe samples in μ g/wipe, soil/solid/sludge samples in mg/kg, product/oil/non-aqueous liquid samples in mg/L, and all DISTLC / STLC / SPLP / TCLP extracts are reported in μ g/L.

cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract.

%SS = Percent Recovery of Surrogate Standard. DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

e1) unmodified or weakly modified diesel is significant

e2) diesel range compounds are significant; no recognizable pattern

Angela Rydelius, Lab Manager

QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57719 WorkOrder 1104507

EPA Method SW8015B Extraction SW3550B/3630C									piked San	nple ID:	: 1104453-0)09A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	1
, and y to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	ND	40	119	121	2.06	93.7	94	0.340	70 - 130	30	70 - 130	30
%SS:	105	25	101	103	1.38	99	99	0	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57719 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104507-001A	04/18/11 12:35 PM	04/18/11	04/22/11 3:51 PM	1104507-002A	04/18/11 12:40 PM	04/18/11	04/19/11 10:56 PM
1104507-003A	04/18/11 12:45 PM	04/18/11	04/20/11 2:19 AM	1104507-004A	04/18/11 1:35 PM	04/18/11	04/20/11 3:27 AM
1104507-005A	04/18/11 1:30 PM	04/18/11	04/19/11 6:24 PM	1104507-006A	04/18/11 4:50 PM	04/18/11	04/19/11 7:32 PM
1104507-007A	04/18/11 4:45 PM	04/18/11	04/19/11 9:48 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

DHS ELAP Certification 1644

QA/QC Officer

QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57739 WorkOrder 1104507

EPA Method SW8021B/8015Bm	Extrac	tion SW	5030B					5	Spiked Sar	nple ID	: 1104507-0	07A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acc	eptance	Criteria (%)	
Allalyto	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex)	ND	0.60	121	118	2.15	124	122	1.91	70 - 130	20	70 - 130	20
MTBE	ND	0.10	113	113	0	113	105	7.33	70 - 130	20	70 - 130	20
Benzene	ND	0.10	91.1	94.1	3.23	92.8	90.4	2.65	70 - 130	20	70 - 130	20
Toluene	ND	0.10	89.2	92	3.04	90.2	88.3	2.09	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	90.5	94.3	4.19	91.4	89.6	2.00	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	90.5	93.7	3.47	91.5	89.5	2.19	70 - 130	20	70 - 130	20
%SS:	83	0.10	89	80	10.0	91	89	2.17	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57739 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104507-001A	04/18/11 12:35 PM	04/18/11	04/20/11 12:55 AM	1104507-002A	04/18/11 12:40 PM	04/18/11	04/20/11 1:25 AM
1104507-003A	04/18/11 12:45 PM	04/18/11	04/20/11 2:54 AM	1104507-004A	04/18/11 1:35 PM	04/18/11	04/20/11 3:23 AM
1104507-005A	04/18/11 1:30 PM	04/18/11	04/20/11 3:53 AM	1104507-006A	04/18/11 4:50 PM	04/18/11	04/20/11 4:22 AM
1104507-007A	04/18/11 4:45 PM	04/18/11	04/20/11 4:52 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

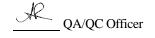
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



QC SUMMARY REPORT FOR SW8270C

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57743 WorkOrder 1104507

EPA Method SW8270C	Extra	ction SW	3550B					S	Spiked Sar	nple ID	: 1104506-0)07A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acc	eptance	Criteria (%))
Analyte	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Acenaphthene	ND	2	70.7	69	2.51	69.6	68.4	1.81	30 - 130	30	30 - 130	30
4-Chloro-3-methylphenol	ND	4	80.2	82.2	2.45	75.1	75.1	0	30 - 130	30	30 - 130	30
2-Chlorophenol	ND	4	70.8	67.9	4.17	75.4	70.2	7.10	30 - 130	30	30 - 130	30
1,4-Dichlorobenzene	ND	2	76.7	73.5	4.14	74.6	74.7	0.0804	30 - 130	30	30 - 130	30
2,4-Dinitrotoluene	ND	2	83.1	82.1	1.17	85	80.2	5.74	30 - 130	30	30 - 130	30
4-Nitrophenol	ND	4	72	71.3	1.05	66.3	69.8	5.17	30 - 130	30	30 - 130	30
N-Nitrosodi-n-propylamine	ND	2	71.8	69.2	3.68	91	79.5	13.5	30 - 130	30	30 - 130	30
Pentachlorophenol	ND	4	66.3	64.5	2.78	60.6	58	4.41	30 - 130	30	30 - 130	30
Phenol	ND	4	69.9	68.5	2.08	79.4	73.1	8.23	30 - 130	30	30 - 130	30
Pyrene	ND	2	88.6	82.2	7.45	87.2	83.6	4.19	30 - 130	30	30 - 130	30
1,2,4-Trichlorobenzene	ND	2	83	81.1	2.34	77.8	79.6	2.29	30 - 130	30	30 - 130	30
%SS1:	94	200	88	85	3.77	101	88	13.9	30 - 130	30	30 - 130	30
%SS2:	105	200	90	87	3.58	106	94	12.1	30 - 130	30	30 - 130	30
%SS3:	96	200	96	96	0	99	96	3.61	30 - 130	30	30 - 130	30
%SS4:	83	200	76	80	5.10	77	76	0.310	30 - 130	30	30 - 130	30
%SS5:	111	200	105	95	10.1	99	97	2.83	30 - 130	30	30 - 130	30
%SS6:	91	200	91	85	7.23	88	92	5.04	30 - 130	30	30 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57743 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104507-001A	04/18/11 12:35 PM	04/18/11	04/21/11 8:03 PM	1104507-002A	04/18/11 12:40 PM	04/18/11	04/21/11 9:20 PM
1104507-003A	04/18/11 12:45 PM	04/18/11	04/21/11 10:36 PM	1104507-004A	04/18/11 1:35 PM	04/18/11	04/21/11 11:52 PM
1104507-005A	04/18/11 1:30 PM	04/18/11	04/22/11 1:07 AM	1104507-006A	04/18/11 4:50 PM	04/18/11	04/22/11 2:22 AM
1104507-007A	04/18/11 4:45 PM	04/18/11	04/21/11 10:16 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and / or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix, sample diluted due to high matrix or analyte content, or MS/MSD samples diluted due to high organic content.

#) surrogate diluted out of range; & = low or no recovery of surrogate or target analytes due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.

QA/QC Officer

QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104507

EPA Method SW6010B		Extraction SW3050B				BatchID: 57701			iked Sample ID: 1104507-007A				
Analyte	Sample	Spiked	MS	MSD	MS-MSD	Spiked	LCS	LCSD	LCS-LCSD	Acc	eptanc	e Criteria (%	o)
Analyto	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Lead	9.3	50	107	108	1.07	10	84.2	92.5	9.43	75 - 125	25	75 - 125	25
%SS:	113	500	113	115	2.28	500	111	114	2.31	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57701 SUMMARY

Lab ID	Date Sampled	Date Extract	ed Date Analyzed	Lab ID	Date Sampled	Date Extracte	d Date Analyzed
1104507-001A	04/18/11 12:35 PM	04/18/11	04/19/11 11:49 AM	1104507-002A	04/18/11 12:40 PM	04/18/11	04/19/11 11:52 AM
1104507-003A	04/18/11 12:45 PM	04/18/11	04/19/11 3:05 PM	1104507-004A	04/18/11 1:35 PM	04/18/11	04/19/11 12:00 PM
1104507-005A	04/18/11 1:30 PM	04/18/11	04/19/11 12:02 PM	1104507-006A	04/18/11 4:50 PM	04/18/11	04/19/11 12:04 PM
1104507-007A	04/18/11 4:45 PM	04/18/11	04/19/11 11:16 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer

McCampbell Analytical,	Inc.
"When Quality Counts"	

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled: 04/20/11
555 Montgomery St., Suite 1300		Date Received: 04/20/11
tee mongomery on, same 1999	Client Contact: Peter Cusack	Date Reported: 04/22/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/22/11

WorkOrder: 1104581

April 22, 2011

D	D .
I lear	Peter:

Enclosed within are:

- 1) The results of the 1 analyzed sample from your project: #730482302; 5812 Hollis St.,
- 2) A QC report for the above sample,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.



Treadwell & Rollo CHAIN OF CUSTODY RECORD

Page ____ of ____

Environmental and Geot	echnical Consult	tant	555 M 501 1																			415.	955.	9041		
Site Name:	5812	Holl	777 C												CA	9582	5 Ph	916	.565.	7412		910	6.565			
Job Number: Project Manager\Co	ntact:	Peter	Cusack		-							(Su)		П	A	naly	sis	Req	ues	ted	Γ			Tui	rnaround Time	
Samplers: Recorder (Signature	Required):	Milan	Tw							tain		2.8	S									ean-up		12	thour	
Field Sample Identification No.	Date	Time	Lab Sample No.	Soil	Water			$\overline{}$	90		Other	TPH-	50/00	ead								Silica gel clean-up	Hold	Remar	ka	
EP-1	4-20-11	1100		X			7	7	7	X	+	X	X	X	1	+	+		1	1		Χ̈́		Nemai	na .	
				,	-			1	1	-	+				1	+	+			+						-
							1	1	1	+	+				+	+	+		+	+						-
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Relinguished by: (Signatu	ire))	Date 4/200/11				Time	19	3	0		Red	eive	d by	: (Sig	gnatu	ire)	,				Date	е	Time		_
Relinquished by: (Signatu			Date				Time	1				Rec	eive	STAY	tab	(Sig	natur	(e) _	2,	7	5	Date	94	20/11 Time /	1430	
Sent to Laboratory (f aboratory Commen		McC	amppell								-	Me				ipme rried		Priva	La ite Co			Na:		ed Ex Airborn	e UPS	
	1	White Copy	· Original		Yello	ow (Сор	y - I	Lab	orate	ory]		HE.	OD C AD S CHL	CONI PACI ORIN	EABS	IN L	AB_	_	CON	ROP TAI SER	RIA' NER	IN LAB	244	

1534 Willow Pass Rd

CHAIN-OF-CUSTODY RECORD

Page 1 of 1

1104581-001	EP-1		Soil	4/20/2011 11:00		Α	Α	Α								
Lab ID	Client ID		Matrix	Collection Date	Hold	1	2	3	4	5	6	7 8	9	10	11	12
									Requ	ested Te	sts (Se	e legend	below)			
-	Rollo omery St., Suite 1300 sco, CA 94111	cc: PO:	,	adwellrollo.com 5812 Hollis St.			Tre 55:	eadwell 5 Montg	Payable & Rollo Jomery S isco, CA	St., Suite	1300	Da	quested te Rece te Prin	ived:		
		WaterTrax	WriteOn	EDF		Excel	l	Fax	✓] Email		HardCopy	Thir	dParty	J-	flag
	rg, CA 94565-1701 52-9262					Work	Order:	11045	581	Cli	entCode	e: TWRF				

Test Legend:				
1 8270D_S	2 PB_S	3 TPH(DMO)WSG_S	4	5
6	7	8	9	10
11	12			
The following SampID: 001A contains	testgroup.			Prepared by: Maria Venegas

Comments: SEND HARD COPY. 72hr Rush

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

Sample Receipt Checklist

Client Name:	Treadwell & Roll	0			Date	and Time Received:	4/20/2011	2:48:12 PM
Project Name:	#730482302; 581	2 Hollis St.			Chec	klist completed and	reviewed by:	Maria Venegas
WorkOrder N°:	1104581	Matrix Soil			Carrie	er: Rob Pringle (N	MAI Courier)	
		<u>Cha</u>	in of Cu	ıstody (C	COC) Informa	ation		
Chain of custody	/ present?		Yes	V	No 🗆			
•	· / signed when relinqui	ished and received?	Yes	V	No 🗆			
	/ agrees with sample l		Yes	✓	No 🗌			
Sample IDs noted	d by Client on COC?		Yes	V	No 🗆			
Date and Time of	f collection noted by Cl	ient on COC?	Yes	~	No 🗆			
Sampler's name r	noted on COC?		Yes	✓	No 🗆			
			Sample	Receip	t Information	n		
Custody seals int	tact on shipping conta		Yes		No 🗆	_	NA 🔽	
-	er/cooler in good cond		Yes	V	No 🗆			
	er containers/bottles?		Yes	~	No 🗆			
Sample containe	ers intact?		Yes	✓	No 🗆			
Sufficient sample	e volume for indicated	test?	Yes	✓	No 🗆			
		Sample Pres	ervatio	n and Ho	old Time (HT) Information		
All samples recei	ived within holding tim	ne?	Yes	✓	No 🗌			
Container/Temp I	Blank temperature		Coole	er Temp:	7.6°C		NA 🗆	
Water - VOA vial	ls have zero headspa	ice / no bubbles?	Yes		No 🗆	No VOA vials subr	nitted 🗹	
Sample labels ch	necked for correct pre	servation?	Yes	~	No 🗌			
Metal - pH accep	otable upon receipt (pF	H<2)?	Yes		No 🗆		NA 🗹	
Samples Receive	ed on Ice?		Yes	✓	No 🗆			
		(Ice Ty	rpe: WE	ET ICE)			
* NOTE: If the "N	No" box is checked, s	ee comments below						
=====		=====	===		====	=====	====	
Client contacted:		Date conta	cted:			Contacted	d by:	
Comments:								

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/20/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received: 04/20/11
555 Wongomery St., State 1500	Client Contact: Peter Cusack	Date Extracted: 04/20/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/21/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104581

Lab ID				1104581-001A			
Client ID				EP-1			
Matrix				Soil			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33
2.4.6-Trichlorophenol	ND	1.0	0.33				
	-	Surre	ogate Re	coveries (%)	-	·	·

Surrogate Recoveries (%)										
%SS1:	89	%SS2:	90							
%SS3:	94	%SS4:	76							
%SS5:	72	%SS6:	91							

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor



^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

McCampbell Analytical, Inc. "When Ouality Counts'

1534 Willow Pass Road, Pittsburg, CA 94565-1701 $Web: www.mccampbell.com \qquad E-mail: main@mccampbell.com\\$ Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo 555 Montgomery St., Suite 1300	Client Project ID: #730482302; 5812 Hollis St.	Date Sampled: 04/20/11
	Hollis St.	Date Received: 04/20/11
	Client Contact: Peter Cusack	Date Extracted: 04/20/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 04/21/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline*

Extraction method SW5	straction method SW5030B		ethods SW8015Bm	Wo	ork Order:	1104581
Lab ID	Client ID	Matrix	TPH(g)	DF	% SS	Comments
001A	EP-1	S	ND	1	86	
	ing Limit for DF =1;	W	NA		NA	
	ans not detected at or the reporting limit	S	1.0		mg/Kg	3

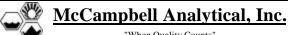
Reporting Limit for DF =1;	W	NA	NA		
ND means not detected at or	S	1.0	mg/Kg		
above the reporting limit	5	1.0	mg/Kg		

^{*} water and vapor samples are reported in µg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts in mg/L.

Angela Rydelius, Lab Manager

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:



"When Ouality Counts"			Te	lephone: 8	377-252-9262 Fax: 925	-252-9	269			
Treadwell & F	Rollo	Client Project ID: Hollis St.		730482302; 5812	2	Date Sampled:	04/2	20/11		
555 Montgomery St., Suite 1300		Hollis St.			Date Received:	04/2	20/11			
		Client Contact: Peter Cusack			Date Extracted:	04/2	20/11			
San Francisco	o, CA 94111	Client P.O.:			Date Analyzed:	04/2	21/11			
Lead by ICP*										
Extraction method	: SW3050B		Analyt	tical methods: SW60)10B				Work Ord	er: 1104581
Lab ID	Client ID		Matrix	Extraction Type		Lead	j	DF	% SS	Comments
1104581-001A	EP-1		S	TOTAL		8.6		1	93	

Reporting Limit for DF =1;	W	TOTAL	NA	μg/L
ND means not detected at or	S	TOTAL	5.0	mg/Kg
above the reporting limit				8 8

*water samples are reported in µg/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, filter samples in µg/filter.

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of $0.45 \mu m$ filtered and acidified sample.

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

Angela Rydelius, Lab Manager

DHS ELAP Certification 1644



	Client Project ID: #730482302; 5812	Date Sampled:	04/20/11
555 Montgomery St., Suite 1300	Hollis St.	Date Received:	04/20/11
	Client Contact: Peter Cusack	Date Extracted:	04/20/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/21/11

Total Extractable Petroleum Hydrocarbons with Silica Gel Clean-Up* Extraction method: SW3550B/3630C Analytical methods: SW8015B Work Order: 1104581 TPH-Diesel TPH-Motor Oil DF % SS Lab ID Client ID Matrix Comments (C10-C23) (C18-C36) 1104581-001A EP-1 S ND ND 105 Reporting Limit for DF =1; W NA NA ug/L ND means not detected at or S 1.0 5.0 mg/Kgabove the reporting limit

* water samples are reported in μ g/L, wipe samples in μ g/wipe, soil/solid/sludge samples in mg/kg, product/oil/non-aqueous liquid samples in mg/L, and all DISTLC / STLC / SPLP / TCLP extracts are reported in μ g/L.

cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract.

%SS = Percent Recovery of Surrogate Standard. DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

DHS ELAP Certification 1644

Angela Rydelius, Lab Manager

QC SUMMARY REPORT FOR SW8270C

QC Matrix: Soil BatchID: 57743 WorkOrder 1104581 W.O. Sample Matrix: Soil

EPA Method SW8270C	Extra	ction SW	3550B					s	Spiked San	nple ID:	1104506-0	07A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
, may to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Acenaphthene	ND	2	70.7	69	2.51	69.6	68.4	1.81	30 - 130	30	30 - 130	30
4-Chloro-3-methylphenol	ND	4	80.2	82.2	2.45	75.1	75.1	0	30 - 130	30	30 - 130	30
2-Chlorophenol	ND	4	70.8	67.9	4.17	75.4	70.2	7.10	30 - 130	30	30 - 130	30
1,4-Dichlorobenzene	ND	2	76.7	73.5	4.14	74.6	74.7	0.0804	30 - 130	30	30 - 130	30
2,4-Dinitrotoluene	ND	2	83.1	82.1	1.17	85	80.2	5.74	30 - 130	30	30 - 130	30
4-Nitrophenol	ND	4	72	71.3	1.05	66.3	69.8	5.17	30 - 130	30	30 - 130	30
N-Nitrosodi-n-propylamine	ND	2	71.8	69.2	3.68	91	79.5	13.5	30 - 130	30	30 - 130	30
Pentachlorophenol	ND	4	66.3	64.5	2.78	60.6	58	4.41	30 - 130	30	30 - 130	30
Phenol	ND	4	69.9	68.5	2.08	79.4	73.1	8.23	30 - 130	30	30 - 130	30
Pyrene	ND	2	88.6	82.2	7.45	87.2	83.6	4.19	30 - 130	30	30 - 130	30
1,2,4-Trichlorobenzene	ND	2	83	81.1	2.34	77.8	79.6	2.29	30 - 130	30	30 - 130	30
%SS1:	94	200	88	85	3.77	101	88	13.9	30 - 130	30	30 - 130	30
%SS2:	105	200	90	87	3.58	106	94	12.1	30 - 130	30	30 - 130	30
%SS3:	96	200	96	96	0	99	96	3.61	30 - 130	30	30 - 130	30
%SS4:	83	200	76	80	5.10	77	76	0.310	30 - 130	30	30 - 130	30
%SS5:	111	200	105	95	10.1	99	97	2.83	30 - 130	30	30 - 130	30
%SS6:	91	200	91	85	7.23	88	92	5.04	30 - 130	30	30 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57743 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed	
1104581-001A	04/20/11 11:00 AM	04/20/11	04/21/11 5:16 PM					

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and / or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix, sample diluted due to high matrix or analyte content, or MS/MSD samples diluted due to high organic content.

#) surrogate diluted out of range; & = low or no recovery of surrogate or target analytes due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels

A QA/QC Officer

QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57720 WorkOrder 1104581

EPA Method SW8015Bm	S	Spiked Sample ID: 1104453-009A										
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
7 mary to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex [£]	ND	0.60	99.5	103	3.37	101	99.6	0.934	70 - 130	20	70 - 130	20
MTBE	ND	0.10	109	106	3.39	106	107	0.515	70 - 130	20	70 - 130	20
Benzene	ND	0.10	105	104	0.374	105	104	1.09	70 - 130	20	70 - 130	20
Toluene	ND	0.10	92.1	92.1	0	92	90.8	1.26	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	94.1	94.6	0.490	93.6	92.4	1.25	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	106	107	0.969	106	104	1.99	70 - 130	20	70 - 130	20
%SS:	97	0.10	108	106	2.74	105	106	1.03	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57720 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104581-001A	04/20/11 11:00 AM	1 04/20/11	04/21/11 3:11 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

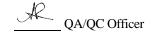
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104581

EPA Method SW6010B			Extract	ion SW3	3050B		BatchID	: 57805	Spiked Sample ID: 1104581-001A							
Analyte	Sample	Spiked	d MS MSD MS-MS		MS-MSD	Spiked	LCS	LCSD	LCS-LCSD	Acc	eptanc	e Criteria (%	5)			
7a.y to	mg/Kg mg/Kg % Rec. % R		% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD				
Lead	8.6	50	89	88.8	0.236	10	98.3	92.8	5.78	75 - 125	25	75 - 125	25			
%SS:	93	500	106	103	2.01	500	102	98	3.25	70 - 130	20	70 - 130	20			

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57805 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104581-001A	04/20/11 11:00 AM	04/20/11	04/21/11 1:44 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer

QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57810 WorkOrder 1104581

EPA Method SW8015B			5	Spiked Sample ID: 1104581-001A								
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	١
, many to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	ND	40	114	114	0	92.6	90.9	1.89	70 - 130	30	70 - 130	30
%SS:	105	25	104	102	1.08	92	93	1.15	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57810 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104581-001A	04/20/11 11:00 AM	04/20/11	04/21/11 1:56 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

DHS ELAP Certification 1644

QA/QC Officer

McCampbell Analytical,	Inc.
"When Quality Counts"	

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled: 04/26/11
555 Montgomery St., Suite 1300		Date Received: 04/26/11
oss frongomery bu, butter 1900	Client Contact: Peter Cusack	Date Reported: 04/27/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/27/11

WorkOrder: 1104725

April 27, 2011

D	D .
I lear	Peter:

Enclosed within are:

- 1) The results of the 1 analyzed sample from your project: #730482302; 5812 Hollis St,
- 2) A QC report for the above sample,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.





Treadwell&Rollo

CHAIN OF CUSTODY RECORD

555 Montgomery Street, Suite 1300, San Francisco, CA 94111 Ph: 415.955.9040/Fax: 415.955.9041

Page __ of __

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Site Name: Job Number:	5812	Hollis 18238	SF		15 0	Omn	nons	Ru	., oui	16 200,	6		T		nal							ax.	910	1]		Turnaro		
Project Manager\Cor Samplers: Recorder (Signature	Rob	Ma	Corace TWO	М	atri	×				iners vative	10 - OL - Cai	20	20	7									gel clean-up			2	4 Time	001	
Field Sample Identification No.	Date	Time	Lab Sample No.	_	Water	$\overline{}$	_	$\overline{}$	_	Other	TICK LY	X	AAT	160									Silica gel o	Hold		Rem	arks		
12 - (-) /	4-26-11	1055		×										7									_						
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<i>*</i>	^					+	+		+		+				D	CH	ER	RIN	ON	VC	LA AS	B_ 0&	-		ESERVED IN L	4B	-		_
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Relinquished by: (Signature) Date						-	Time						ed b	10	_	_0	7	1	/	1	4	1	Date	4	126/11	Time	170	_	
Sent to Laboratory (Laboratory Commen	nts/Notes:	Mc	anpoell								M	ethe	od o	nd C	arrie	d [ate	Cou	cou			me)		Airbo		UPS	3
		White Copy	- Original		Yell	ow	Copy	y - L	abo	ratory					Pinl	CC	ру	- F	ielo	1			C	000	C Number:	00!	524!	5	

McCampbell Analytical, Inc.

Report to:

1534 Willow Pass Rd Pittsburg, CA 94565-1701 (925) 252-9262

CHAIN-OF-CUSTODY RECORD

Page 1 of 1

WorkOrder: 1104725 ClientCode: TWRF

HardCopy

WaterTrax	WriteOn	☐ EDF	Excel	Fax	✓ Email	HardCopy ThirdParty	J-flag
			Bill	to:		Requested TAT:	1 day

Peter Cusack Email: pjcusack@treadwellrollo.com Treadwell & Rollo cc:

□ WaterTrax

555 Montgomery St., Suite 1300 PO:

San Francisco, CA 94111 ProjectNo: #730482302; 5812 Hollis St

FAX (415) 955-9041 (415) 955-5244

Accounts Payable Treadwell & Rollo

Fax

555 Montgomery St., Suite 1300

San Francisco, CA 94111

Date Received: 04/26/2011

Date Printed: 04/26/2011

				Requested Tests (See legend below)											
Lab ID	Client ID	Matrix	Collection Date Hold	1	2	3	4	5	6	7	8	9	10	11	12
1104725-001	B-1-17	Soil	4/26/2011 10:55	Α	Α	Α									

Test Legend:

1	G-MBTEX_S	2	PB_S	3 TPH(DMO)WSG_S]	4	5
6		7		8]	9	10
11		12					
							Prepared by: Ana Venegas

Comments: SEND HARD COPY 24hr rush

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

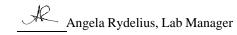
Sample Receipt Checklist

Client Name: Treadwell & Rollo				Date ar	nd Time Received:	4/26/2011	5:30:40 PM			
Project Name: #730482302; 5812 Holl	lis St			Checkl	ist completed and re	eviewed by:	Ana Venegas			
WorkOrder N°: 1104725 Matrix	C Soil			Carrier	: Rob Pringle (M.	AI Courier)				
	<u>Chain o</u>	of Cu	stody (C	OC) Informat	<u>tion</u>					
Chain of custody present?	,	Yes	V	No 🗆						
Chain of custody signed when relinquished a	nd received?	Yes	V	No 🗆						
Chain of custody agrees with sample labels?	,	Yes	✓	No 🗌						
Sample IDs noted by Client on COC?		Yes	V	No 🗆						
Date and Time of collection noted by Client on	COC?	Yes	~	No 🗆						
Sampler's name noted on COC?	,	Yes	V	No 🗆						
	<u>Sar</u>	nple	Receipt	<u>Information</u>						
Custody seals intact on shipping container/co		Yes		No 🗆		NA 🔽				
Shipping container/cooler in good condition?	,	Yes	V	No 🗆						
Samples in proper containers/bottles?		Yes	✓	No 🗆						
Sample containers intact?		Yes	✓	No 🗆						
Sufficient sample volume for indicated test?	,	Yes	✓	No 🗌						
Sample Preservation and Hold Time (HT) Information										
All samples received within holding time?	,	Yes	✓	No 🗌						
Container/Temp Blank temperature	(Coole	er Temp:	7°C		NA 🗆				
Water - VOA vials have zero headspace / no	bubbles?	Yes		No 🗆	No VOA vials submi	tted 🗹				
Sample labels checked for correct preservati	on?	Yes	✓	No 🗌						
Metal - pH acceptable upon receipt (pH<2)?	,	Yes		No 🗆		NA 🗹				
Samples Received on Ice?		Yes	✓	No 🗆						
	(Ice Type:	WE	TICE))						
* NOTE: If the "No" box is checked, see com	ments below.									
	=====	==	===							
Client contacted:	Date contacted	d:			Contacted	by:				
Comments:										

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled:	04/26/11
555 Montgomery St., Suite 1300	Hollis St	Date Received:	04/26/11
• •	Client Contact: Peter Cusack	Date Extracted:	04/26/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/27/11

	Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline with BTEX and MTBE*										
Extraction	on method: SW5030B			Analy	tical methods:	SW8021B/8015	Bm		Wor	k Order:	1104725
Lab ID	Client ID	Matrix	TPH(g)	MTBE	Benzene	Toluene	Ethylbenzene	Xylenes	DF	% SS	Comments
001A	B-1-17	S	ND	ND	ND	ND	ND	ND	1	86	
	ting Limit for DF =1;	W	50	5.0	0.5	0.5	0.5	0.5		ug/I	,
	eans not detected at or re the reporting limit	S	1.0	0.05	0.005	0.005	0.005	0.005		mg/K	

* water and vapor samples are reported in µg/L, soil/sludge/solid	samples in mg/kg,	wipe samples in µg/wipe,	product/oil/non-aqueous	liquid samples a	and all
TCLP & SPLP extracts in mg/L.					



[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:



Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled: 04/26/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/26/11
<i>y</i> ,	Client Contact: Peter Cusack	Date Extracted: 04/26/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/27/11

Lead by ICP*

Extraction method: SW3050B Analytical methods: SW6010B Work Order: 1104725

Extraction method.	3 W 2020D	Anarytear methods. Sw0010B						
Lab ID	Client ID	Matrix	Extraction Type	Lead	DF	% SS	Comments	
1104725-001A	B-1-17	S	TOTAL	5.2	1	108		

Reporting Limit for DF =1;	W	TOTAL	NA	μg/L
ND means not detected at or	S	TOTAL	5.0	mg/Kg
above the reporting limit	5	TOTAL	3.0	mg Kg

*water samples are reported in µg/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, filter samples in µg/filter.

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of $0.45 \mu m$ filtered and acidified sample.

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

DHS ELAP Certification 1644

Angela Rydelius, Lab Manager



1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

		Date Sampled:	04/26/11
555 Montgomery St., Suite 1300	Hollis St	Date Received:	04/26/11
	Client Contact: Peter Cusack	Date Extracted:	04/26/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/27/11

San Francisco, CA	A 94111	Client P.O.:		Date Analyzed:	04/27/	11	
Extraction method: SW			m Hydrocarbons with	h Silica Gel Clean-Up*	W	ork Order:	1104725
Lab ID	Client ID	Matrix	TPH-Diesel (C10-C23)	TPH-Motor Oil (C18-C36)	DF	% SS	Comments
1104725-001A	B-1-17	S	ND	ND	1	101	
	ing Limit for DF =1;	W	NA	NA		ug/.	L
ND mea	ans not detected at or	S	1.0	5.0		ma/l	Vα

above the reporting limit	3	1.0	5.0	mg/Kg
* water samples are reported in µg/L, wipe samples in µg.	wipe, soil/	solid/sludge samples in mg/	kg, product/oil/non-aqueous	liquid samples in mg/L, and all
DISTLC / STLC / SPLP / TCLP extracts are reported in	$\mu g/L$.			

1.0

cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract.

%SS = Percent Recovery of Surrogate Standard. DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

DHS ELAP Certification 1644

Angela Rydelius, Lab Manager

5.0

mg/Kg

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

QC SUMMARY REPORT FOR SW8021B/8015Bm

QC Matrix: Soil BatchID: 57866 WorkOrder 1104725 W.O. Sample Matrix: Soil

EPA Method SW8021B/8015Bm	Extra	tion SW	5030B			Spiked Sample ID: 1104662-004A												
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	CSD Acceptance Criteria (%)									
Allalyte	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD LCS/LCS		RPD						
TPH(btexf	ND	0.60	92.1	99.3	7.57	93.3	84.2	10.2	70 - 130	20	70 - 130	20						
MTBE	ND	0.10	95.2	94.7	0.485	100	86.1	15.0	70 - 130	20	70 - 130	20						
Benzene	ND	0.10	89.8	94.8	5.43	101	92.4	8.62	70 - 130	20	70 - 130	20						
Toluene	ND	0.10	79.8	86.1	7.59	88.7	82.4	7.34	70 - 130	20	70 - 130	20						
Ethylbenzene	ND	0.10	82.4	89.9	8.65	90.4	84.4	6.87	70 - 130	20	70 - 130	20						
Xylenes	ND	0.30	94.4	104	10.0	102	95.9	6.21	70 - 130	20	70 - 130	20						
%SS:	92	0.10	90	102	12.5	108	99	9.45	70 - 130	20	70 - 130	20						

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57866 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed	
1104725-001A	04/26/11 10:55 AM	1 04/26/11	04/27/11 11:32 AM					

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

A QA/QC Officer

QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104725

EPA Method SW6010B			Extract	ion SW3	3050B		BatchID	: 57934	Spik	ed Sample	ID:	1104725-00	1A
Analyte	Sample	Spiked	MS	MS MSD MS		Spiked	LCS	LCSD	LCS-LCSD	Acc	eptanc	e Criteria (%	5)
Amaryto	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Lead	5.2	50	112	102	8.96	10	101	101	0	75 - 125	25	75 - 125	25
%SS:	108	500	113	110	2.42	500	108	109	1.52	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57934 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104725-001A	04/26/11 10:55 AM	04/26/11	04/27/11 1:34 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57933 WorkOrder 1104725

EPA Method SW8015B	Extrac	tion SW	3550B/36	630C				S									
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD									
, many to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD					
TPH-Diesel (C10-C23)	ND	40	125	126	1.09	104	97.5	6.21	70 - 130	30	70 - 130	30					
%SS:	101	25	117	117	0	93	86	7.53			30						

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57933 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed	
1104725-001A	04/26/11 10:55 AM	I 04/26/11	04/27/11 9:01 AM					

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

DHS ELAP Certification 1644

QA/QC Officer

McCampbell Analytical,	Inc.
"When Quality Counts"	

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled: 04/28/11
555 Montgomery St., Suite 1300		Date Received: 04/28/11
oss frongomery bu, butter 1900	Client Contact: Peter Cusack	Date Reported: 04/29/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/29/11

WorkOrder: 1104805

April 29, 2011

D	D .
I lear	Peter:

Enclosed within are:

- 1) The results of the 1 analyzed sample from your project: #730482302; 5812 Hollis St,
- 2) A QC report for the above sample,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.





Treadwe Environmental and Geote			CHAI	ontgo 4th S	omer	y Str	reet	, Suit	e 13 Oal	300, Sakland	an F	Fran 946	cisc 12 F	o, C	A 94	74.4	Ph: 500/	Fax	: 51	0.8	74.	4507	7							Page	10	if
Site Name: Job Number:	5812	182302	777 Ca	ampu -	is Co	mmo	ons	Rd.,	Suit	te 200	, Sa	acra	men	ito, (alys							x: 9	16.	.565.	7412		-	Tu	rnaro	und	
Project Manager∖Cor Samplers: Recorder (Signature	P	PCUSAL VISALL	work	м	atrix					iners		dyno	9	4									done in	dn-uear			4		2	Time	OVE	
Field Sample Identification No.	Date	Time	Lab Sample No.	_	$\overline{}$	Other	_	HNO	lce	Other	\neg	A 149	201/8	154									Olive on it	Silica gei clean-up	Hold			R	emar	ks		
13-15-17.51	4/28/11	11:2ZA	4	X		+	+		X		1	4	4	1	+	-	F		F	t	+	+	+	+	+							
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Relinguished by: (Signatu Relinguisheoloy: (Signatu	an)	Date 4/29/11				ime	13	1/2	5	1		agencies in	-	(Sig	forting.			-	 ×	2		1	ate	1/0	28/	//	4	Time	13	15	_
Relinquished by: (Signatu	ure)		Date /			П	ime	10			F	Regi	eivet	by	Lab:	(Sign	natu	re)	2	j			D	ate	4	128,	10	,	Time	143	20	
Sent to Laboratory (Laboratory Commen		McCi	tay hell							_	-	Met	_		Shi _l Car			Pri	_	_		ouri		Van		ed Ex			Airborn	е [U	PS
		White Copy	- Original		Yello	ow C	Ор	y - La	abo	ratory	/				P	ink (Сор	y - F	Fie	d				C	ОС	Numl	ber:	0	05	226	6	

McCampbell Analytical, Inc.

1534 Willow Pass Rd Pittsburg, CA 94565-1 (925) 252-9262

CHAIN-OF-CUSTODY RECORD

Page 1 of 1

	g, CA 94565-1701 52-9262					Work	Order	11048	805	C	lientCo	ode: TV	VRF				
		WaterTrax	WriteOn	☐ EDF		Excel	[Fax	5	✓ Email		HardC	ору	Thire	dParty	□J-f	lag
Report to: Peter Cusad	ck	Email: p	ojcusack@tre	adwellrollo.com		١	Bill to: Ac	counts	Payable	Э			Requ	ıested	TAT:	1	day
•	mery St., Suite 1300	cc: PO:					55	5 Montg		St., Suit					- ,	04/28/	
San Francis (415) 955-52	co, CA 94111 44 FAX (415) 955-9041	ProjectNo: #	±730482302;	5812 Hollis St			Sa	n Franc	cisco, C	A 94111			Date	Print	ted:	04/28/	2011
									Requ	uested 7	Гests (See lege	end be	elow)			
Lab ID	Client ID		Matrix	Collection Date	Hold	1	2	3	4	5	6	7	8	9	10	11	12
1104805-001	B-15-17.5'		Soil	4/28/2011 11:22		Α	Α	А									

Test Legend:				
1 8270D_S	2 G-MBTEX_S	3 PB_S	4	5
6	7	8	9	10
11	12			
The following SampID: 001A cont	ains testgroup.		_	Prepared by: Ana Venegas

Comments: SEND HARD COPY 24hr rush

NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days).

Hazardous samples will be returned to client or disposed of at client expense.

Sample Receipt Checklist

Client Name: Treadwell &	Rollo			Date ar	nd Time Received:	4/28/2011	2:25:36 PM
Project Name: #730482302;	5812 Hollis St			Checkl	ist completed and rev	riewed by:	Ana Venegas
WorkOrder N°: 1104805	Matrix Soil			Carrier	: Rob Pringle (MA	I Courier)	
	<u>Chair</u>	n of Cu	ıstody (C	COC) Informat	<u>tion</u>		
Chain of custody present?		Yes	V	No 🗆			
Chain of custody signed when rel	inquished and received?	Yes	V	No \square			
Chain of custody agrees with sam	nple labels?	Yes	✓	No 🗌			
Sample IDs noted by Client on COO	0?	Yes	V	No 🗆			
Date and Time of collection noted I	by Client on COC?	Yes	~	No 🗆			
Sampler's name noted on COC?		Yes	✓	No 🗆			
	s	ample	Receipt	Information			
Custody seals intact on shipping of		Yes		No 🗆	N	IA 🔽	
Shipping container/cooler in good	condition?	Yes	V	No 🗆			
Samples in proper containers/bott	tles?	Yes	✓	No 🗆			
Sample containers intact?		Yes	✓	No 🗆			
Sufficient sample volume for indic	ated test?	Yes	✓	No 🗌			
	Sample Prese	rvatio	n and Ho	old Time (HT)	<u>Information</u>		
All samples received within holdin	g time?	Yes	✓	No 🗆			
Container/Temp Blank temperature	e	Coole	er Temp:	10.2°C	١	va 🗆	
Water - VOA vials have zero hea	dspace / no bubbles?	Yes		No 🗆	No VOA vials submitt	ed 🗹	
Sample labels checked for correct	t preservation?	Yes	✓	No 🗌			
Metal - pH acceptable upon receip	ot (pH<2)?	Yes		No \square	١	va 🔽	
Samples Received on Ice?		Yes	~	No 🗆			
	(Ice Typ	e: WE	ET ICE)			
* NOTE: If the "No" box is checke	ed, see comments below.						
========	=======	=					
Client contacted:	Date contac	ted:			Contacted b	y:	
Comments:							

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/28/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/28/11
333 Wongomery St., State 1300	Client Contact: Peter Cusack	Date Extracted: 04/28/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 04/28/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1104805

Lab ID	1104805-001A								
Client ID		B-15-17.5'							
Matrix		Soil							
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit		
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33		
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33		
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6		
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33		
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33		
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6		
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33		
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33		
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33		
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66		
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33		
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33		
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33		
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33		
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33		
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66		
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33		
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33		
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6		
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33		
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33		
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33		
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33		
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33		
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33		
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33		
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33		
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6		
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33		
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6		
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33		
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33		
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33		
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33		
2.4.6-Trichlorophenol	ND	1.0	0.33						
				coveries (%)					

	Surrogate R	ecoveries (%)	
%SS1:	100	%SS2:	105
%SS3:	109	%SS4:	94
%SS5:	91	%SS6:	81
	·	·	

Comments:

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.

^{*} water samples in $\mu g/L$, soil/sludge/solid samples in mg/kg, wipe samples in $\mu g/wipe$, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: #730482302; 5812	Date Sampled: 04/28/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/28/11
555 Hongomery Bu, Butte 1500	Client Contact: Peter Cusack	Date Extracted: 04/28/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 04/29/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline*

Extraction method	SW5030B		al methods SW8015Bm	Wo	ork Order:	1104805
Lab ID	Client ID	Matrix	TPH(g)	DF	% SS	Comments
001A	B-15-17.5'	S	79	10	76	ď7
	porting Limit for DF =1;	W	NA		NA	
	means not detected at or bove the reporting limit	S	1.0		mg/Kg	3

^{*} water and vapor samples are reported in µg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts in mg/L.

d7) strongly aged gasoline or diesel range compounds are significant in the TPH(g) chromatogram

Angela Rydelius, Lab Manager

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:



	•		
Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled:	04/28/11
555 Montgomery St., Suite 1300	Hollis St	Date Received:	04/28/11
	Client Contact: Peter Cusack	Date Extracted:	04/28/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/28/11
	Lead by ICP*		
Extraction method: SW3050B	Analytical methods: SW6010B		Work Order: 1104805

Extraction method: SW3	0B Analytical methods: SW6010B						der: 1104805
Lab ID	Client ID	Matrix	Extraction Type	Lead	DF	% SS	Comments
1104805-001A	B-15-17.5'	S	TOTAL	9.1	1	109	

Reporting Limit for DF =1;	W	TOTAL	NA	μg/L
ND means not detected at or	S	TOTAL	5.0	mg/Kg
above the reporting limit	~	TOTALE	3.0	mg ng

*water samples are reported in µg/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, filter samples in µg/filter.

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of $0.45 \mu m$ filtered and acidified sample.

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

Angela Rydelius, Lab Manager

DHS ELAP Certification 1644

McCampbell Analytical, Inc.
"When Ouality Counts"

	Client Project ID: #730482302; 5812	Date Sampled:	04/28/11
555 Montgomery St., Suite 1300	Hollis St	Date Received:	04/28/11
	Client Contact: Peter Cusack	Date Extracted:	04/28/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	04/28/11

Extraction method: SW35	50B		nethods: SW8015B	ocarbons*	Work Order: 1104805				
Lab ID	Client ID	Matrix	TPH-Diesel (C10-C23)	TPH-Motor Oil (C18-C36)	DF	% SS	Comments		
1104805-001A	B-15-17.5'	S	270	61	1	110	e1		
Reporting	g Limit for DF =1;	W	NA	NA		ug/l	 Г.		

Reporting Limit for DF =1;	W	NA	NA	ug/L
ND means not detected at or above the reporting limit	S	1.0	5.0	mg/Kg

^{*} water samples are reported in μ g/L, wipe samples in μ g/wipe, soil/solid/sludge samples in mg/kg, product/oil/non-aqueous liquid samples in mg/L, and all DISTLC / STLC / SPLP / TCLP extracts are reported in μ g/L.

cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

e1) unmodified or weakly modified diesel is significant

Angela Rydelius, Lab Manager

DHS ELAP Certification 1644

QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57866 WorkOrder 1104805

EPA Method SW8015Bm	EPA Method SW8015Bm Extraction SW5030B Spiked Sample ID: 1104662-004A											
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acc	eptance	Criteria (%)	
7 mary to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex)	ND	0.60	92.1	99.3	7.57	93.3	84.2	10.2	70 - 130	20	70 - 130	20
MTBE	ND	0.10	95.2	94.7	0.485	100	86.1	15.0	70 - 130	20	70 - 130	20
Benzene	ND	0.10	89.8	94.8	5.43	101	92.4	8.62	70 - 130	20	70 - 130	20
Toluene	ND	0.10	79.8	86.1	7.59	88.7	82.4	7.34	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	82.4	89.9	8.65	90.4	84.4	6.87	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	94.4	104	10.0	102	95.9	6.21	70 - 130	20	70 - 130	20
%SS:	92	0.10	90	102	12.5	108	99	9.45	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57866 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104805-001A	04/28/11 11:22 AM	1 04/28/11	04/29/11 11:16 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

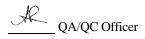
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



QC SUMMARY REPORT FOR SW8270C

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57948 WorkOrder 1104805

EPA Method SW8270C	EPA Method SW8270C Extraction SW3550B Spiked Sample ID: 1104750-017A)17A		
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acceptance Criteria (%)				
, mary to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD	
Acenaphthene	ND<1.6	2	71.2	71.2	0	75.2	76.3	1.42	30 - 130	30	30 - 130	30	
4-Chloro-3-methylphenol	ND<1.6	4	75.8	76	0.198	77.6	74.8	3.74	30 - 130	30	30 - 130	30	
2-Chlorophenol	ND<1.6	4	89.2	93.3	4.58	92.5	91.5	1.09	30 - 130	30	30 - 130	30	
1,4-Dichlorobenzene	ND<1.6	2	92.2	93.6	1.61	91.3	90.2	1.17	30 - 130	30	30 - 130	30	
2,4-Dinitrotoluene	ND<1.6	2	77.8	76.8	1.29	87.6	85.6	2.37	30 - 130	30	30 - 130	30	
4-Nitrophenol	ND<8.0	4	93.3	92.4	0.996	84.8	76.2	10.6	30 - 130	30	30 - 130	30	
N-Nitrosodi-n-propylamine	ND<1.6	2	111	111	0	109	107	1.79	30 - 130	30	30 - 130	30	
Pentachlorophenol	ND<8.0	4	61.6	56.2	9.12	32.2	34.9	8.19	30 - 130	30	30 - 130	30	
Phenol	ND<1.6	4	65	63.6	2.18	85.8	85.6	0.140	30 - 130	30	30 - 130	30	
Pyrene	ND<1.6	2	71.9	71.6	0.488	75	77.1	2.68	30 - 130	30	30 - 130	30	
1,2,4-Trichlorobenzene	ND<1.6	2	83.8	85.2	1.72	88.2	87.8	0.546	30 - 130	30	30 - 130	30	
%SS1:	82	200	82	83	1.06	84	82	2.47	30 - 130	30	30 - 130	30	
%SS2:	81	200	69	71	1.91	90	90	0	30 - 130	30	30 - 130	30	
%SS3:	89	200	89	90	0.877	98	98	0	30 - 130	30	30 - 130	30	
%SS4:	78	200	79	79	0	87	87	0	30 - 130	30	30 - 130	30	
%SS5:	95	200	100	103	2.13	102	101	0.286	30 - 130	30	30 - 130	30	
%SS6:	72	200	72	72	0	80	80	0	30 - 130	30	30 - 130	30	

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57948 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104805-001A	04/28/11 11:22 AM	1 04/28/11	04/28/11 5:49 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

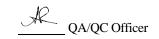
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and / or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix, sample diluted due to high matrix or analyte content, or MS/MSD samples diluted due to high organic content.

#) surrogate diluted out of range; & = low or no recovery of surrogate or target analytes due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.



QC SUMMARY REPORT FOR 6010B

W.O. Sample Matrix: Soil QC Matrix: Soil WorkOrder 1104805

EPA Method SW6010B Extraction SW3050B				3050B	BatchID: 57934 Spiked Sample ID: 1104725-00					1A			
Analyte		Spiked	MS	MSD	MS-MSD	Spiked	LCS	LCSD	LCS-LCSD	Acc	eptanc	e Criteria (%	5)
, many to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	mg/Kg	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Lead	5.2	50	112	102	8.96	10	101	101	0	75 - 125	25	75 - 125	25
%SS:	108	500	113	110	2.42	500	108	109	1.52	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57934 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104805-001A	04/28/11 11:22 AM	04/28/11 (4/28/11 10:17 PM			•	

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer

QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57932 WorkOrder 1104805

EPA Method SW8015B	3550B					S	piked San	nple ID:	: 1104723-0	002A		
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%))
, many to	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	6.0	40	106	108	1.39	106	104	2.07	70 - 130	30	70 - 130	30
%SS:	100	25	103	103	0	99	94	5.30	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57932 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed	
1104805-001A	04/28/11 11:22 AM	I 04/28/11	04/28/11 6:55 PM					

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

DHS ELAP Certification 1644

QA/QC Officer

McCampbell Ar "When Quality		1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269				
Treadwell & Rollo	Client Project ID: 5812 Holl	is St	Date Sampled:	05/03/11		
555 Montgomery St., Suite 1300			Date Received:	05/03/11		
555 Honegomery Su, Suite 1500	Client Contact: Peter Cusa	ck	Date Reported:	05/04/11		
San Francisco, CA 94111	Client P.O.:		Date Completed:	05/04/11		

WorkOrder: 1105078

May 04, 2011

Dear P	eter:
--------	-------

Enclosed within are:

- 1) The results of the 1 analyzed sample from your project: 5812 Hollis St,
- 2) A QC report for the above sample,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.

1105078



White Copy - Original

CHAIN OF CUSTODY RECORD

Page ___ of __

Job Number:		Hollis													Ar	aly	/sis	Re	qu	iest	ed						Turnaround
Project Manager\Con Samplers:	5000 - 000					_						MO												dr			24 110UR
Recorder (Signature	Required):								ont			9	2							1				clean-up		L	1
			where the same		latr	$\overline{}$	$\overline{}$	$\overline{}$	esei	$\overline{}$	$\overline{}$	Ha.	R											gel c			
Field Sample Identification No.	Date.	Time	Lab Sample No.	Soil	Water	Othe	된	H ₂ SO ₄	N -	eo loe	e co	4	51											Silica	Hold		Remarks
B-15-19	5/3/11	11:00 40		Ϋ́,				4		4	\perp	X	X			1	\perp	L	1	\perp	1	1	-	X			
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				+		Н	\dashv	+	+	+	+	\vdash		Н	\vdash	+	+	+	+	+	$^{+}$	+	+	\dashv	_		
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4/0	/										\perp						1	1			1	1				//	
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Retinquished by Signat	ire)		Date / 3///				Tim	7	3	>		Red	ceive	ed by	y: (&i	matu	ure)				\	V	9	Date	e/	/	Time
Relinquished by: (Signatu	ure)	11 0	Date / //				Tim	9				Red	ceive	d by	y ab	: (Sig	gnatu	ire)	1	/		4	10	Date	e <	3/11	TIME 730
Sent to Laboratory (Name): nts/Notes:	Mac	AMPHS									Me	tho	1	f Shi	-		_	X	La	ab c	couri	ier	[me)	Fed Ex	Airborne UPS

Yellow Copy - Laboratory

Pink Copy - Field

COC Number: 005227

McCampbell Analytical, Inc.

1534 Willow Pass Rd

CHAIN-OF-CUSTODY RECORD

Page 1 of 1

Pittsburg, CA 94565-1701 (925) 252-9262					Work	Orde	er: 110	5078	3	C	ClientC	ode: T	WRF				
	WaterTrax	WriteOn	☐ EDF		Excel		Fax	<	✓	Email		Hard	Сору	Thir	dParty	□J-	flag
Report to:						Bill to	o:						Req	uested	TAT:	1	day
Peter Cusack Treadwell & Rollo 555 Montgomery St., Suite 1300 San Francisco, CA 94111 (415) 955-5244 FAX (415) 955-9041	cc: PO: ProjectNo: {	ojcusack@tre 5812 Hollis Si	adwellrollo.com			7 5	Account Freadwe 555 Mor San Frai	ell & ntgom	Rollo nery S)		e Rece e Prini		05/03/ 05/03/	
									Requ	ested	Tests (See le	gend b	elow)	,		
Lab ID Client ID		Matrix	Collection Date	Hold	1	2	3		4	5	6	7	8	9	10	11	12
1105078-001 B-15-9		Soil	5/3/2011 11:00		Α	Α											
Test Legend:																	

The following SampID: 001A contains testgroup.

8270D_S

SEND HARD COPY 24hr rush **Comments:**

2

7

12

TPH(DMO)WSG_S

NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

3

8

5

Prepared by: Ana Venegas

Sample Receipt Checklist

Client Name:	Treadwell & R	ollo			Date a	and Time Received:	5/3/2011 7	7:41:29 PM
Project Name:	5812 Hollis St				Check	klist completed and r	eviewed by:	Ana Venegas
WorkOrder N°:	1105078	Matrix <u>Soil</u>			Carrie	er: Rob Pringle (M	IAI Courier)	
		<u>Cha</u>	in of Cu	ıstody (C	COC) Informa	ation_		
Chain of custody	y present?		Yes	V	No 🗆			
Chain of custody	y signed when relin	quished and received?	Yes	V	No \square			
Chain of custody	y agrees with samp	le labels?	Yes	✓	No 🗌			
Sample IDs noted	d by Client on COC?		Yes	V	No 🗆			
Date and Time of	f collection noted by	Client on COC?	Yes	V	No 🗆			
Sampler's name	noted on COC?		Yes		No 🔽			
			Sample	Receip	t Information	1		
Custody seals in	ntact on shipping co		Yes		No 🗆	-	NA 🔽	
Shipping contain	ner/cooler in good co	ondition?	Yes	V	No 🗆			
Samples in prop	er containers/bottle	s?	Yes	V	No 🗆			
Sample containe	ers intact?		Yes	✓	No 🗆			
Sufficient sample	e volume for indicat	ed test?	Yes	✓	No 🗌			
		Sample Pres	ervatio	n and H	old Time (HT) Information		
All samples rece	eived within holding	-	Yes	V	No 🗆	<u>, </u>		
	Blank temperature		Coole	er Temp:	5.2°C		NA 🗆	
		pace / no bubbles?	Yes		No 🗆	No VOA vials subm	itted 🗹	
	hecked for correct		Yes	~	No 🗌			
Metal - pH accep	otable upon receipt	(pH<2)?	Yes		No \square		NA 🗹	
Samples Receive	ed on Ice?		Yes	V	No \square			
		(Ice Ty	/pe: WE	ET ICE)			
* NOTE: If the "I	No" box is checked	, see comments below						
=====	=====	======	===	===	====	======		======
Client contacted:	:	Date conta	cted:			Contacted	by:	
Comments:								

Treadwell & Rollo	Client Project ID: 5812 Hollis St	Date Sampled: 05/03/11
555 Montgomery St., Suite 1300		Date Received: 05/03/11
333 Montgomery St., Suite 1300	Client Contact: Peter Cusack	Date Extracted: 05/03/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 05/04/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3550B Analytical Method: SW8270C Work Order: 1105078

Lab ID		1105078-001A									
Client ID				B-15-19							
Matrix				Soil							
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit				
Acenaphthene	ND	1.0	0.33	Acenaphthylene	ND	1.0	0.33				
Acetochlor	ND	1.0	0.33	Anthracene	ND	1.0	0.33				
Benzidine	ND	1.0	1.6	Benzoic Acid	ND	1.0	1.6				
Benzo(a)anthracene	ND	1.0	0.33	Benzo(b)fluoranthene	ND	1.0	0.33				
Benzo(k)fluoranthene	ND	1.0	0.33	Benzo(g,h,i)perylene	ND	1.0	0.33				
Benzo(a)pyrene	ND	1.0	0.33	Benzyl Alcohol	ND	1.0	1.6				
1,1-Biphenyl	ND	1.0	0.33	Bis (2-chloroethoxy) Methane	ND	1.0	0.33				
Bis (2-chloroethyl) Ether	ND	1.0	0.33	Bis (2-chloroisopropyl) Ether	ND	1.0	0.33				
Bis (2-ethylhexyl) Phthalate	ND	1.0	0.33	4-Bromophenyl Phenyl Ether	ND	1.0	0.33				
Butylbenzyl Phthalate	ND	1.0	0.33	4-Chloroaniline	ND	1.0	0.66				
4-Chloro-3-methylphenol	ND	1.0	0.33	2-Chloronaphthalene	ND	1.0	0.33				
2-Chlorophenol	ND	1.0	0.33	4-Chlorophenyl Phenyl Ether	ND	1.0	0.33				
Chrysene	ND	1.0	0.33	Dibenzo(a,h)anthracene	ND	1.0	0.33				
Dibenzofuran	ND	1.0	0.33	Di-n-butyl Phthalate	ND	1.0	0.33				
1,2-Dichlorobenzene	ND	1.0	0.33	1,3-Dichlorobenzene	ND	1.0	0.33				
1,4-Dichlorobenzene	ND	1.0	0.33	3,3-Dichlorobenzidine	ND	1.0	0.66				
2,4-Dichlorophenol	ND	1.0	0.33	Diethyl Phthalate	ND	1.0	0.33				
2,4-Dimethylphenol	ND	1.0	0.33	Dimethyl Phthalate	ND	1.0	0.33				
4,6-Dinitro-2-methylphenol	ND	1.0	1.6	2,4-Dinitrophenol	ND	1.0	1.6				
2,4-Dinitrotoluene	ND	1.0	0.33	2,6-Dinitrotoluene	ND	1.0	0.33				
Di-n-octyl Phthalate	ND	1.0	0.33	1,2-Diphenylhydrazine	ND	1.0	0.33				
Fluoranthene	ND	1.0	0.33	Fluorene	ND	1.0	0.33				
Hexachlorobenzene	ND	1.0	0.33	Hexachlorobutadiene	ND	1.0	0.33				
Hexachlorocyclopentadiene	ND	1.0	1.6	Hexachloroethane	ND	1.0	0.33				
Indeno (1,2,3-cd) pyrene	ND	1.0	0.33	Isophorone	ND	1.0	0.33				
2-Methylnaphthalene	ND	1.0	0.33	2-Methylphenol (o-Cresol)	ND	1.0	0.33				
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	0.33	Naphthalene	ND	1.0	0.33				
2-Nitroaniline	ND	1.0	1.6	3-Nitroaniline	ND	1.0	1.6				
4-Nitroaniline	ND	1.0	1.6	Nitrobenzene	ND	1.0	0.33				
2-Nitrophenol	ND	1.0	1.6	4-Nitrophenol	ND	1.0	1.6				
N-Nitrosodiphenylamine	ND	1.0	0.33	N-Nitrosodi-n-propylamine	ND	1.0	0.33				
Pentachlorophenol	ND	1.0	1.6	Phenanthrene	ND	1.0	0.33				
Phenol	ND	1.0	0.33	Pyrene	ND	1.0	0.33				
1,2,4-Trichlorobenzene	ND	1.0	0.33	2,4,5-Trichlorophenol	ND	1.0	0.33				
2.4.6-Trichlorophenol	ND	1.0	0.33								
		Surre	gate Re	coveries (%)							

%SS5: 67

%SS2

%SS4:

80

83

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.

72

69

%SS1:

%SS3:

^{*} water samples in μ g/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

Treadwell & Rollo	Client Project ID: 5812 Hollis St	Date Sampled: 05/03/11
555 Montgomery St., Suite 1300		Date Received: 05/03/11
, ,	Client Contact: Peter Cusack	Date Extracted: 05/03/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 05/04/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline *

	Gasoline Ra	nge (C6-C12) Vo	olatile Hydrocarbons as Gasoline [:]	*		
extraction method SW50)30B	Analytic	al methods SW8015Bm	Wo	ork Order:	1105078
Lab ID	Client ID	Matrix	TPH(g)	DF	% SS	Commen
001A	B-15-19	S	27	1	101	d7
	ng Limit for DF =1; ns not detected at or	W	NA		NA	
	the reporting limit	S	1.0		mg/K	g

above the reporting limit	3	1.0	mg/Kg				
* water and vapor samples are reported in µg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous l							
samples and all TCLP & SPLP extracts in mg/L.							

d7) strongly aged gasoline or diesel range compounds are significant in the TPH(g) chromatogram

Angela Rydelius, Lab Manager

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:



Treadwell & Rollo	Client Project ID: 5812 Hollis St	Date Sampled:	05/03/11
555 Montgomery St., Suite 1300		Date Received:	05/03/11
	Client Contact: Peter Cusack	Date Extracted:	05/03/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed:	05/04/01

San Francisco, C	CA 94111	Client P.O.: Date Analyzed		Date Analyzed:	05/04/	01	
Extraction method: S			m Hydrocarbons wit	h Silica Gel Clean-Up*	W	ork Order:	1105078
Lab ID	Client ID	Matrix	TPH-Diesel (C10-C23)	TPH-Motor Oil (C18-C36)	DF	% SS	Comments
1105078-001A	B-15-19	S	100	31	1	108	e1
Repo	rting Limit for DF =1;	W	NA	NA		ug/.	I

* water samples are reported in	$\mu g/L$, wipe samples in μg	/wipe, soil	/solid/sludge samples in mg/	kg, product/oil/non-aqueous	liquid samples in mg/L, and all
DISTLC / STLC / SPLP / TCL	P extracts are reported in	цg/L.			

1.0

cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract.

%SS = Percent Recovery of Surrogate Standard. DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

e1) unmodified or weakly modified diesel is significant

ND means not detected at or

above the reporting limit

Angela Rydelius, Lab Manager

5.0

DHS ELAP Certification 1644

mg/Kg

QC SUMMARY REPORT FOR SW8270C

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 58035 WorkOrder 1105078

EPA Method SW8270C	Extraction SW3550B						Spiked Sample ID: N/A					
Analyte	Sample	Sample Spiked MS			MS-MSD	LCS	LCSD	LCS-LCSD	Acceptance Criteria (%)			
Analyte	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Acenaphthene	N/A	2	N/A	N/A	N/A	67.5	68.2	1.11	N/A	N/A	30 - 130	30
4-Chloro-3-methylphenol	N/A	4	N/A	N/A	N/A	75.1	77	2.56	N/A	N/A	30 - 130	30
2-Chlorophenol	N/A	4	N/A	N/A	N/A	78.2	80.5	2.93	N/A	N/A	30 - 130	30
1,4-Dichlorobenzene	N/A	2	N/A	N/A	N/A	76.4	78.6	2.92	N/A	N/A	30 - 130	30
2,4-Dinitrotoluene	N/A	2	N/A	N/A	N/A	67.6	67.9	0.398	N/A	N/A	30 - 130	30
4-Nitrophenol	N/A	4	N/A	N/A	N/A	50.2	49.9	0.579	N/A	N/A	30 - 130	30
N-Nitrosodi-n-propylamine	N/A	2	N/A	N/A	N/A	60	63.4	5.48	N/A	N/A	30 - 130	30
Pentachlorophenol	N/A	4	N/A	N/A	N/A	30	30.3	0.995	N/A	N/A	30 - 130	30
Phenol	N/A	4	N/A	N/A	N/A	70.3	72.8	3.56	N/A	N/A	30 - 130	30
Pyrene	N/A	2	N/A	N/A	N/A	85.8	76	12.1	N/A	N/A	30 - 130	30
1,2,4-Trichlorobenzene	N/A	2	N/A	N/A	N/A	85.4	86.7	1.58	N/A	N/A	30 - 130	30
%SS1:	N/A	200	N/A	N/A	N/A	72	71	2.37	N/A	N/A	30 - 130	30
%SS2:	N/A	200	N/A	N/A	N/A	71	73	3.43	N/A	N/A	30 - 130	30
%SS3:	N/A	200	N/A	N/A	N/A	77	81	4.79	N/A	N/A	30 - 130	30
%SS4:	N/A	200	N/A	N/A	N/A	67	67	0	N/A	N/A	30 - 130	30
%SS5:	N/A	200	N/A	N/A	N/A	62	63	0.392	N/A	N/A	30 - 130	30
%SS6:	N/A	200	N/A	N/A	N/A	77	74	4.70	N/A	N/A	30 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 58035 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1105078-001A	05/03/11 11:00 AM	1 05/03/11	05/04/11 2:42 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

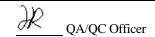
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and / or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix, sample diluted due to high matrix or analyte content, or MS/MSD samples diluted due to high organic content.

#) surrogate diluted out of range; & = low or no recovery of surrogate or target analytes due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.



QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57996 WorkOrder 1105078

EPA Method SW8015Bm	Extraction SW5030B						Spiked Sample ID: 1104808-002A					
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS LCSD		LCS-LCSD	S-LCSD Acceptance Criteria			1
	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex)	ND	0.60	121	118	2.51	117	113	3.66	70 - 130	20	70 - 130	20
MTBE	ND	0.10	98.2	102	4.16	93.4	106	12.6	70 - 130	20	70 - 130	20
Benzene	ND	0.10	89.2	92.1	3.27	88.5	95.5	7.55	70 - 130	20	70 - 130	20
Toluene	ND	0.10	88.8	91.3	2.74	87.8	94	6.85	70 - 130	20	70 - 130	20
Ethylbenzene	ND	0.10	89.7	92.8	3.49	88.8	95.2	7.04	70 - 130	20	70 - 130	20
Xylenes	ND	0.30	89.7	92.3	2.86	88	94.4	7.00	70 - 130	20	70 - 130	20
%SS:	90	0.10	79	81	2.75	80	84	5.70	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57996 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1105078-001A	05/03/11 11:00 AM	I 05/03/11	05/04/11 7:16 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

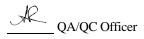
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Soil QC Matrix: Soil BatchID: 57986 WorkOrder 1105078

EPA Method SW8015B	Extraction SW3550B/3630C				Spiked Sample ID: 1104794-026A							
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acceptance Criteria (%)			١
	mg/Kg	mg/Kg	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	ND	40	106	106	0	100	103	2.69	70 - 130	30	70 - 130	30
%SS:	109	25	91	90	0.783	84	88	4.51	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57986 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed	
1105078-001A	05/03/11 11:00 AM	1 05/03/11	05/04/01 3:07 PM					

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

DHS ELAP Certification 1644

QA/QC Officer

McCampbell Analytical, Inc.

1534 Willow Pass Road, Pittsburg, CA 94565-1701
Web: www.mccampbell.com E-mail: main@mccampbell.com
Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; Alders Greenway	Date Sampled: 02/08/11
555 Montgomery St., Suite 1300		Date Received: 02/10/11
tee mongomery on, same 1999	Client Contact: Peter Cusack	Date Reported: 02/11/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 02/11/11

WorkOrder: 1102303

February 11, 2011

D	D .
I lear	Peter:

Enclosed within are:

- 1) The results of the 1 analyzed sample from your project: #730482302; Alders Greenway,
- 2) A QC report for the above sample,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.

RUSH

1102303

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Site Name:	Aldens	6 ROUN	777.0	4th S	stre	et, T	hird	Flo	or, C	Daklar	nd C	A 94	612	Ph:	510	.874	.450	0/F	ax:	51	0.8	74.	4507	7			35.7412	77 50 507			
Job Number:	730 4	82302	,												Α	nal	ysi	s R	lec	que	est	ed						Т	urnaro	und	
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Relinquished by Signat	ure)		Date 2/10/1)	,		ľ	Time	Ki	2	1		Reg	eive	ed by	(Si	gnat	ure)	4	4					D	ate	1	10/11	Time	30		
Relinquished by: (Signat	ure)	11	Date //				Time					Rec	eive	ed by	/ Lat	: (Si	gnat	ure	3					D	ate	, ,		Time			
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McCampbell Analytical, Inc.

1534 Willow Pass Rd Pittsburg, CA 94565-1701 (925) 252-9262

CHAIN-OF-CUSTODY RECORD

ClientCode: TWRF

WorkOrder: 1102303

Page 1 of 1

		WaterTrax	WriteOn	☐ EDF		Excel		Fax	[✓ Email		HardCo	ру	Third	dParty	J-1	flag
Report to:						E	Bill to:						Requ	uested 7	TAT:	1	day
-	Rollo omery St., Suite 1300 sco, CA 94111	cc: PO: ProjectNo: #		adwellrollo.com Alders Greenway			Tre 558 Sa	•	& Rolle gomery cisco, C	o St., Sui A 9411		,		e Recei e Printo		02/10/2 02/10/2	
									Req	uested	Tests (See lege	nd b	elow)			
Lab ID	Client ID		Matrix	Collection Date	Hold	1	2	3	Req 4	uested 5	Tests (See lege 7	nd bo	elow)	10	11	12
Lab ID 1102303-001	Client ID DW-1		Matrix Water	Collection Date	Hold	1 D	2 B	3	Req 4	1		See lege 7			10	11	12

Test Legend:

1	8260B_W		2 8270D_W	3	CAM17MS_DISS	4	G-MBTEX_W		5	PRDISSOLVED
6			7	8		9			10	
11		1	2							
Th	e following SampID: 001A con	tains testo	iroun					Pren	ared l	hv• Zoraida Cortez

Comments:

Sample Receipt Checklist

Client Name:	Treadwell & Rollo				Date a	nd Time Received:	2/10/2011	3:29:59 PM
Project Name:	#730482302; Alders	Greenway			Check	list completed and r	eviewed by:	Zoraida Cortez
WorkOrder N°:	1102303 Mat	rix <u>Water</u>			Carrie	r: Rob Pringle (M	(1Al Courier)	
		<u>Chain</u>	of Cu	stody (C	COC) Informa	<u>tion</u>		
Chain of custody	y present?		Yes	V	No 🗆			
Chain of custody	y signed when relinquished	and received?	Yes	V	No 🗆			
Chain of custody	y agrees with sample labels	s?	Yes	✓	No 🗌			
Sample IDs note	d by Client on COC?		Yes	V	No 🗆			
Date and Time o	of collection noted by Client o	n COC?	Yes	~	No 🗆			
Sampler's name	noted on COC?		Yes	✓	No 🗆			
		<u>S</u>	ample	Receipt	Information			
Custody seals in	ntact on shipping container/o	cooler?	Yes		No 🗆		NA 🔽	
Shipping contain	ner/cooler in good condition?	?	Yes	V	No 🗆			
Samples in prop	per containers/bottles?		Yes		No 🗹			
Sample containe	ers intact?		Yes	✓	No 🗆			
Sufficient sample	e volume for indicated test?		Yes	✓	No 🗌			
		Sample Prese	rvatio	n and Ho	old Time (HT)	Information		
All samples rece	eived within holding time?		Yes	✓	No 🗌			
Container/Temp	Blank temperature		Coole	er Temp:	4.8°C		NA \square	
Water - VOA via	als have zero headspace / r	no bubbles?	Yes		No \square	No VOA vials subm	nitted 🗹	
Sample labels c	hecked for correct preserva	ition?	Yes	✓	No 🗌			
Metal - pH accep	ptable upon receipt (pH<2)?		Yes		No 🗆		NA 🔽	
Samples Receiv	red on Ice?		Yes	✓	No 🗆			
		(Ice Typ	e: WE	TICE)			
* NOTE: If the "	No" box is checked, see co	mments below.						
	=======							
Client contacted	:	Date contact	ed:			Contacted	l by:	
Comments:	No Voas were received.							

Treadwell & Rollo	Client Project ID: #730482302; Alders	Date Sampled: 02/08/11
555 Montgomery St., Suite 1300	Greenway	Date Received: 02/10/11
	Client Contact: Peter Cusack	Date Extracted: 02/10/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 02/10/11

Volatile Organics by P&T and GC/MS (Basic Target List)*

Extraction Method: SW5030B Analytical Method: SW8260B Work Order: 1102303

Extraction Method: SW5030B		Anaiyi	icai Metiic	0d: SW8200B	work Order: 1102	2303	
Lab ID				1102303-001D			
Client ID				DW-1			
Matrix				Water			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acetone	120	1.0	10	tert-Amyl methyl ether (TAME)	ND	1.0	0.5
Benzene	ND	1.0	0.5	Bromobenzene	ND	1.0	0.5
Bromochloromethane	ND	1.0	0.5	Bromodichloromethane	ND	1.0	0.5
Bromoform	ND	1.0	0.5	Bromomethane	ND	1.0	0.5
2-Butanone (MEK)	79	1.0	2.0	t-Butyl alcohol (TBA)	ND	1.0	2.0
n-Butyl benzene	ND	1.0	0.5	sec-Butyl benzene	ND	1.0	0.5
tert-Butyl benzene	ND	1.0	0.5	Carbon Disulfide	0.85	1.0	0.5
Carbon Tetrachloride	ND	1.0	0.5	Chlorobenzene	ND	1.0	0.5
Chloroethane	ND	1.0	0.5	Chloroform	ND	1.0	0.5
Chloromethane	ND	1.0	0.5	2-Chlorotoluene	ND	1.0	0.5
4-Chlorotoluene	ND	1.0	0.5	Dibromochloromethane	ND	1.0	0.5
1.2-Dibromo-3-chloropropane	ND	1.0	0.2	1.2-Dibromoethane (EDB)	ND	1.0	0.5
Dibromomethane	ND	1.0	0.5	1,2-Dichlorobenzene	ND	1.0	0.5
1.3-Dichlorobenzene	ND	1.0	0.5	1.4-Dichlorobenzene	ND	1.0	0.5
Dichlorodifluoromethane	ND	1.0	0.5	1,1-Dichloroethane	ND	1.0	0.5
1,2-Dichloroethane (1,2-DCA)	ND	1.0	0.5	1.1-Dichloroethene	ND	1.0	0.5
cis-1,2-Dichloroethene	ND	1.0	0.5	trans-1,2-Dichloroethene	ND	1.0	0.5
1,2-Dichloropropane	ND	1.0	0.5	1,3-Dichloropropane	ND	1.0	0.5
2.2-Dichloropropane	ND	1.0	0.5	1.1-Dichloropropene	ND	1.0	0.5
cis-1,3-Dichloropropene	ND	1.0	0.5	trans-1,3-Dichloropropene	ND	1.0	0.5
Diisopropyl ether (DIPE)	ND	1.0	0.5	Ethylbenzene	ND	1.0	0.5
Ethyl tert-butyl ether (ETBE)	ND	1.0	0.5	Freon 113	ND	1.0	10
Hexachlorobutadiene	ND	1.0	0.5	Hexachloroethane	ND	1.0	0.5
2-Hexanone	ND	1.0	0.5	Isopropylbenzene	ND	1.0	0.5
4-Isopropyl toluene	ND	1.0	0.5	Methyl-t-butyl ether (MTBE)	ND	1.0	0.5
Methylene chloride	ND	1.0	0.5	4-Methyl-2-pentanone (MIBK)	ND	1.0	0.5
Naphthalene	ND	1.0	0.5	n-Propyl benzene	ND	1.0	0.5
Styrene	ND	1.0	0.5	1.1.1.2-Tetrachloroethane	ND	1.0	0.5
1,1,2,2-Tetrachloroethane	ND	1.0	0.5	Tetrachloroethene	ND	1.0	0.5
Toluene	1.4	1.0	0.5	1,2,3-Trichlorobenzene	ND	1.0	0.5
1,2,4-Trichlorobenzene	ND	1.0	0.5	1,1,1-Trichloroethane	ND	1.0	0.5
1,1,2-Trichloroethane	ND	1.0	0.5	Trichloroethene	ND	1.0	0.5
Trichlorofluoromethane	ND	1.0	0.5	1,2,3-Trichloropropane	ND	1.0	0.5
1,2,4-Trimethylbenzene	0.81	1.0	0.5	1,3,5-Trimethylbenzene	ND	1.0	0.5
Vinvl Chloride	ND	1.0	0.5	Xvlenes	2.0	1.0	0.5
		Surr	ogate Re	ecoveries (%)			
%SS1:	6		G	%SS2:	10	00	
%SS3:	8			/0552.	1 10	,,,	
Comments: c2	. 0			<u> </u>			

Comments: c2

ND means not detected above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

surrogate diluted out of range or coelutes with another peak; &) low surrogate due to matrix interference.

c2) estimated value due to low surrogate recovery, caused by matrix interference.

^{*} water and vapor samples are reported in $\mu g/L$, soil/sludge/solid samples in mg/kg, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L, wipe samples in $\mu g/kg$.

McCampbell Analytical, Inc.

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	3	Date Sampled: 02/08/11
555 Montgomery St., Suite 1300	Greenway	Date Received: 02/10/11
333 Montgomery St., Suite 1300	Client Contact: Peter Cusack	Date Extracted: 02/10/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed: 02/11/11

Semi-Volatile Organics by GC/MS (Basic Target List)*

Extraction Method: SW3510C Analytical Method: SW8270C Work Order: 1102303

Lab ID	1102303-001B							
Client ID				DW-1				
Matrix				Water				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acenaphthene	ND	1.0	10	Acenaphthylene	ND	1.0	10	
Acetochlor	ND	1.0	10	Anthracene	ND	1.0	10	
Benzidine	ND	1.0	50	Benzoic Acid	ND	1.0	50	
Benzo(a)anthracene	ND	1.0	10	Benzo(b)fluoranthene	ND	1.0	10	
Benzo(k)fluoranthene	ND	1.0	10	Benzo(g,h,i)perylene	ND	1.0	10	
Benzo(a)pyrene	ND	1.0	10	Benzyl Alcohol	ND	1.0	50	
1,1-Biphenyl	ND	1.0	10	Bis (2-chloroethoxy) Methane	ND	1.0	10	
Bis (2-chloroethyl) Ether	ND	1.0	10	Bis (2-chloroisopropyl) Ether	ND	1.0	10	
Bis (2-ethylhexyl) Phthalate	ND	1.0	20	4-Bromophenyl Phenyl Ether	ND	1.0	10	
Butylbenzyl Phthalate	ND	1.0	10	4-Chloroaniline	ND	1.0	20	
4-Chloro-3-methylphenol	ND	1.0	10	2-Chloronaphthalene	ND	1.0	10	
2-Chlorophenol	ND	1.0	10	4-Chlorophenyl Phenyl Ether	ND	1.0	10	
Chrysene	ND	1.0	10	Dibenzo(a,h)anthracene	ND	1.0	10	
Dibenzofuran	ND	1.0	10	Di-n-butyl Phthalate	ND	1.0	10	
1,2-Dichlorobenzene	ND	1.0	10	1,3-Dichlorobenzene	ND	1.0	10	
1,4-Dichlorobenzene	ND	1.0	10	3,3-Dichlorobenzidine	ND	1.0	20	
2,4-Dichlorophenol	ND	1.0	10	Diethyl Phthalate	ND	1.0	10	
2,4-Dimethylphenol	ND	1.0	10	Dimethyl Phthalate	ND	1.0	10	
4,6-Dinitro-2-methylphenol	ND	1.0	50	2,4-Dinitrophenol	ND	1.0	50	
2,4-Dinitrotoluene	ND	1.0	10	2,6-Dinitrotoluene	ND	1.0	10	
Di-n-octyl Phthalate	ND	1.0	10	1,2-Diphenylhydrazine	ND	1.0	10	
Fluoranthene	ND	1.0	10	Fluorene	ND	1.0	10	
Hexachlorobenzene	ND	1.0	10	Hexachlorobutadiene	ND	1.0	10	
Hexachlorocyclopentadiene	ND	1.0	50	Hexachloroethane	ND	1.0	10	
Indeno (1,2,3-cd) pyrene	ND	1.0	10	Isophorone	ND	1.0	10	
2-Methylnaphthalene	ND	1.0	10	2-Methylphenol (o-Cresol)	ND	1.0	10	
3 &/or 4-Methylphenol (m,p-Cres	ND	1.0	10	Naphthalene	ND	1.0	10	
2-Nitroaniline	ND	1.0	50	3-Nitroaniline	ND	1.0	50	
4-Nitroaniline	ND	1.0	50	Nitrobenzene	ND	1.0	10	
2-Nitrophenol	ND	1.0	50	4-Nitrophenol	ND	1.0	50	
N-Nitrosodiphenylamine	ND	1.0	10	N-Nitrosodi-n-propylamine	ND	1.0	10	
Pentachlorophenol	ND	1.0	50	Phenanthrene	ND	1.0	10	
Phenol	29	1.0	10	Pyrene	ND	1.0	10	
1,2,4-Trichlorobenzene	ND	1.0	10	2,4,5-Trichlorophenol	ND	1.0	10	
2.4.6-Trichlorophenol	ND	1.0	10					
		Surro	gate Re	coveries (%)				

Surrogate Recoveries (%)								
%SS1:	94	%SS2:	89					
%SS3:	90	%SS4:	71					
%SS5:	95	%SS6:	92					

ND means not detected at or above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

#) surrogate diluted out of range or surrogate coelutes with another peak.



^{*} water samples in µg/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L.

"When Ouality Counts" Telephone: 877-252-9262 Fax: 925-252-9269 Treadwell & Rollo Client Project ID: #730482302; Alders Date Sampled: 02/08/11 Greenway Date Received: 02/10/11 555 Montgomery St., Suite 1300 Client Contact: Peter Cusack Date Extracted: 02/10/11 San Francisco, CA 94111 Client P.O.: Date Analyzed 02/10/11 CAM / CCR 17 Metals* Lab ID 1102303-001C Reporting Limit for DF =1; ND means not detected Client ID DW-1 above the reporting limit Matrix W S Extraction Type DISS. mg/kg μg/L ICP-MS Metals, Concentration* Work Order: 1102303 Analytical Method: E200.8 Extraction Method: E200.8 Dilution Factor 1 0.5 Antimony ND NA ND 0.5 Arsenic NA 5.0 Barium ND NA Beryllium ND NA 0.5 0.25 ND NA Cadmium 0.5 Chromium ND NA ND NA 0.5 Cobalt ND NA 0.5 Copper Lead ND NA 0.5 0.025Mercury ND NA Molybdenum ND NA 0.5 Nickel ND NA 0.5 0.5 Selenium ND NA 0.19 Silver ND NA 0.5 Thallium ND NA Vanadium 0.5 ND NA 5.0 Zinc ND NA %SS: N/A

*water samples are reported in μ g/L, product/oil/non-aqueous liquid samples and all TCLP / STLC / DISTLC / SPLP extracts are reported in mg/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, filter samples in μ g/filter.

means surrogate diluted out of range; ND means not detected above the reporting limit/method detection limit; N/A means not applicable to this sample or instrument.

TOTAL = Hot acid digestion of a representative sample aliquot.

TRM = Total recoverable metals is the "direct analysis" of a sample aliquot taken from its acid-preserved container.

DISS = Dissolved metals by direct analysis of 0.45 μm filtered and acidified sample.

%SS = Percent Recovery of Surrogate Standard

DF = Dilution Factor

Comments

Treadwell & Rollo	Client Project ID: #730482302; Alders	Date Sampled: 02/08/11
555 Montgomery St., Suite 1300	Greenway	Date Received: 02/10/11
	Client Contact: Peter Cusack	Date Extracted: 02/10/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 02/10/11

San Tancisco	0, 011) +1111	chefit I .O		Date Maryzea 02	/10/11	
	Gasoline Ran	ge (C6-C12) V	olatile Hydrocarbons as Ga	asoline*		_
Extraction method	SW5030B	Analytic	al methods SW8015Bm	Wo	rk Order:	1102303
Lab ID	Client ID	Matrix	TPH(g)	DF	% SS	Comments
001A	DW-1	W	63	1	106	d2
	eporting Limit for DF =1;	W	50		μg/L	
	O means not detected at or above the reporting limit	S	NA		NA	

	· · ·	30	μgL
ND means not detected at or	C	N A	NI A
above the reporting limit	S	INA	NA
*	1/1:41		/-:1/1::4

^{*} water and vapor samples are reported in ug/L, soil/sludge/solid samples in mg/kg, wipe samples in µg/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts in mg/L.

cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference.

%SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

d2) heavier gasoline range compounds are significant (aged gasoline?)

Treadwell & Rollo

Client Project ID: #730482302; Alders
Greenway

Date Sampled: 02/08/11

Date Received: 02/10/11

Client Contact: Peter Cusack

Date Extracted: 02/10/11

San Francisco, CA 94111

Client Project ID: #730482302; Alders
Date Sampled: 02/10/11

Date Received: 02/10/11

Total Extractable Petroleum Hydrocarbons*

Extraction method SW3510C Analytical methods: SW8015B Work Order: 1102303

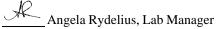
Extraction method 3 v	733100	Anarytica	ai inclious. Sw6013D		WOIK OIU	1. 1102303
Lab ID	Client ID	Matrix	TPH-Diesel (C10-C23)	DF	% SS	Comments
1102303-001A	DW-1	W	97	1	99	e2
	ig Limit for DF =1;	W S	50		με	
	ND means not detected at or above the reporting limit		NA		N	A

^{*} water samples are reported in ug/L, wipe samples in μ g/wipe, soil/solid/sludge samples in mg/kg, product/oil/non-aqueous liquid samples in mg/L, and all DISTLC / STLC / SPLP / TCLP extracts are reported in μ g/L.

%SS = Percent Recovery of Surrogate Standard. DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

e2) diesel range compounds are significant; no recognizable pattern



[#] cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract.

QC SUMMARY REPORT FOR SW8260B

W.O. Sample Matrix: Water QC Matrix: Water BatchID: 56180 WorkOrder 1102303

EPA Method SW8260B	Extra	ction SW	5030B					S	Spiked Sar	nple ID	: 1102304-0	001B
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acc	Criteria (%))	
7 way to	μg/L	μg/L	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
tert-Amyl methyl ether (TAME)	ND<50	10	82.7	84.7	2.32	75.4	82.9	9.49	70 - 130	30	70 - 130	30
Benzene	ND<50	10	105	105	0	96.3	99	2.76	70 - 130	30	70 - 130	30
t-Butyl alcohol (TBA)	ND<200	50	76.4	77.1	0.952	87.7	92.4	5.27	70 - 130	30	70 - 130	30
Chlorobenzene	ND<50	10	105	108	2.65	94.5	104	9.89	70 - 130	30	70 - 130	30
1,2-Dibromoethane (EDB)	ND<50	10	99.6	100	0.739	88.5	95.4	7.50	70 - 130	30	70 - 130	30
1,2-Dichloroethane (1,2-DCA)	ND<50	10	93.7	94.3	0.606	96	99	3.12	70 - 130	30	70 - 130	30
1,1-Dichloroethene	ND<50	10	111	111	0	115	115	0	70 - 130	30	70 - 130	30
Diisopropyl ether (DIPE)	ND<50	10	102	103	1.00	103	106	2.89	70 - 130	30	70 - 130	30
Ethyl tert-butyl ether (ETBE)	ND<50	10	92	93	1.16	97.5	99.7	2.18	70 - 130	30	70 - 130	30
Methyl-t-butyl ether (MTBE)	ND<50	10	98.8	101	2.51	96.6	100	3.48	70 - 130	30	70 - 130	30
Toluene	ND<50	10	100	101	0.222	91.5	95.4	4.13	70 - 130	30	70 - 130	30
Trichloroethene	ND<50	10	105	107	1.76	97.3	99.5	2.25	70 - 130	30	70 - 130	30
%SS1:	96	25	89	89	0	94	93	1.55	70 - 130	30	70 - 130	30
%SS2:	99	25	102	102	0	99	98	0.0989	70 - 130	30	70 - 130	30
%SS3:	77	2.5	83	83	0	86	83	3.37	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 56180 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1102303-001D	02/08/1	1 02/10/11	02/10/11 4:57 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

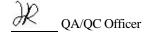
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

surrogate diluted out of range or coelutes with another peak; &) low surrogate due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.



QC SUMMARY REPORT FOR SW8270C

W.O. Sample Matrix: Water QC Matrix: Water BatchID: 56135 WorkOrder 1102303

EPA Method SW8270C Extraction SW3510C Spiked Sample ID: N/A												
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	Acceptance Criteria (%)		
7 mary to	μg/L	μg/L	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Acenaphthene	N/A	50	N/A	N/A	N/A	55.8	59.1	5.76	N/A	N/A	30 - 130	20
4-Chloro-3-methylphenol	N/A	100	N/A	N/A	N/A	82.3	84.9	3.09	N/A	N/A	30 - 130	20
2-Chlorophenol	N/A	100	N/A	N/A	N/A	72.1	75.9	5.15	N/A	N/A	30 - 130	20
1,4-Dichlorobenzene	N/A	50	N/A	N/A	N/A	48.4	51.4	5.94	N/A	N/A	30 - 130	20
2,4-Dinitrotoluene	N/A	50	N/A	N/A	N/A	78.1	80.3	2.83	N/A	N/A	30 - 130	20
4-Nitrophenol	N/A	100	N/A	N/A	N/A	83.7	83.5	0.287	N/A	N/A	30 - 130	20
N-Nitrosodi-n-propylamine	N/A	50	N/A	N/A	N/A	82.5	84.9	2.86	N/A	N/A	30 - 130	20
Pentachlorophenol	N/A	100	N/A	N/A	N/A	68.1	68.9	1.10	N/A	N/A	30 - 130	20
Phenol	N/A	100	N/A	N/A	N/A	78.3	78.8	0.726	N/A	N/A	30 - 130	20
Pyrene	N/A	50	N/A	N/A	N/A	63.4	66.4	4.54	N/A	N/A	30 - 130	20
1,2,4-Trichlorobenzene	N/A	50	N/A	N/A	N/A	54.1	56.3	3.91	N/A	N/A	30 - 130	20
%SS1:	N/A	5000	N/A	N/A	N/A	86	89	3.65	N/A	N/A	30 - 130	20
%SS2:	N/A	5000	N/A	N/A	N/A	80	85	5.76	N/A	N/A	30 - 130	20
%SS3:	N/A	5000	N/A	N/A	N/A	89	93	5.26	N/A	N/A	30 - 130	20
%SS4:	N/A	5000	N/A	N/A	N/A	55	58	5.20	N/A	N/A	30 - 130	20
%SS5:	N/A	5000	N/A	N/A	N/A	96	98	1.97	N/A	N/A	30 - 130	20
%SS6:	N/A	5000	N/A	N/A	N/A	65	67	4.17	N/A	N/A	30 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 56135 SUMMARY

	Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
Γ	1102303-001B	02/08/11	1 02/10/11	02/11/11 4:40 AM			•	

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

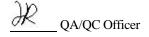
% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.



QC SUMMARY REPORT FOR E200.8

W.O. Sample Matrix: Water QC Matrix: Water BatchID: 56124 WorkOrder 1102303

EPA Method E200.8 Extraction E200.8 Spiked Sample ID: 1102173-003A												
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acc	eptance	Criteria (%)	
7 mary to	μg/L	μg/L	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
Antimony	ND	10	101	105	3.73	98.4	98	0.458	70 - 130	20	85 - 115	20
Arsenic	2.4	10	100	107	5.30	98.9	98.9	0	70 - 130	20	85 - 115	20
Barium	93	100	99.5	107	4.08	97.8	97	0.821	70 - 130	20	85 - 115	20
Beryllium	ND	10	91.7	92.9	1.35	100	98.5	1.65	70 - 130	20	85 - 115	20
Cadmium	ND	10	101	103	1.57	102	101	0.787	70 - 130	20	85 - 115	20
Chromium	ND	10	97	101	4.34	100	100	0	70 - 130	20	85 - 115	20
Cobalt	ND	10	95	96.5	1.56	106	105	1.52	70 - 130	20	85 - 115	20
Copper	91	10	NR	NR	NR	99.6	97.6	1.99	70 - 130	20	85 - 115	20
Lead	ND	10	103	105	2.46	98.5	98.5	0	70 - 130	20	85 - 115	20
Mercury	ND	0.25	94.7	103	8.69	90.2	86.9	3.70	70 - 130	20	85 - 115	20
Molybdenum	3.3	10	104	108	2.67	98.2	97.3	0.920	70 - 130	20	85 - 115	20
Nickel	1.6	10	93.6	99.5	5.26	103	103	0	70 - 130	20	85 - 115	20
Selenium	0.54	10	102	103	0.649	112	107	5.01	70 - 130	20	85 - 115	20
Silver	ND	10	97.4	101	3.31	101	101	0	70 - 130	20	85 - 115	20
Thallium	ND	10	100	104	3.62	96.3	95.8	0.562	70 - 130	20	85 - 115	20
Vanadium	2.8	10	108	113	3.26	98.7	100	1.48	70 - 130	20	85 - 115	20
Zinc	5.4	100	101	105	3.78	107	106	1.12	70 - 130	20	85 - 115	20
%SS:	99	750	99	106	6.99	96	99	2.99	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 56124 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1102303-001C	02/08/11	02/10/11	02/10/11 5:11 PM				

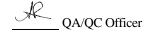
MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not applicable to this method.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Water QC Matrix: Water BatchID: 56186 WorkOrder 1102303

EPA Method SW8015B	Spiked Sample ID: N/A											
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	١
, many to	μg/L	μg/L	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	N/A	1000	N/A	N/A	N/A	114	113	1.34	N/A	N/A	70 - 130	30
%SS:	N/A	625	N/A	N/A	N/A	100	99	0.691	N/A	N/A	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 56186 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1102303-001A	02/08/11	02/10/11	02/10/11 10:51 PM				

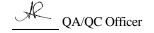
MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.



QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Water QC Matrix: Water BatchID: 56187 WorkOrder 1102303

EPA Method SW8021B/8015Bm Extraction SW5030B Spiked Sample ID: 1102303-001A												
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acceptance Criteria (%)			
7 tildiyto	μg/L	μg/L	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex)	ND	60	97.3	118	19.5	112	98.5	12.7	70 - 130	20	70 - 130	20
MTBE	ND	10	79	81	1.56	122	119	2.09	70 - 130	20	70 - 130	20
Benzene	ND	10	105	107	1.63	108	108	0	70 - 130	20	70 - 130	20
Toluene	1.6	10	92	93.2	1.11	108	108	0	70 - 130	20	70 - 130	20
Ethylbenzene	ND	10	103	105	1.83	106	105	0.103	70 - 130	20	70 - 130	20
Xylenes	3.4	30	98.9	101	2.18	109	109	0	70 - 130	20	70 - 130	20
%SS:	106	10	99	101	2.51	99	98	1.40	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 56187 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1102303-001A	02/08/1	1 02/10/11	02/10/11 8:55 PM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

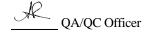
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content, or inconsistency in sample containers.



McCampbell Analytical,	Inc.
"When Ovelity Counts"	

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled: 04/26/11
555 Montgomery St., Suite 1300		Date Received: 04/26/11
tee mongomery on, same 1999	Client Contact: Peter Cusack	Date Reported: 04/27/11
San Francisco, CA 94111	Client P.O.:	Date Completed: 04/27/11

WorkOrder: 1104728

April 27, 2011

D	D .
I lear	Peter:

Enclosed within are:

- 1) The results of the 1 analyzed sample from your project: #730482302; 5812 Hollis St,
- 2) A QC report for the above sample,
- 3) A copy of the chain of custody, and
- 4) An invoice for analytical services.

All analyses were completed satisfactorily and all QC samples were found to be within our control limits.

If you have any questions or concerns, please feel free to give me a call. Thank you for choosing

McCampbell Analytical Laboratories for your analytical needs.

Best regards,

Angela Rydelius Laboratory Manager

McCampbell Analytical, Inc.

1104728

RUSH

CHAIN OF CUSTODY RECORD

Page __ of __

Environmental and Geot	echnical Consult	tant									00, Sa land C														115.	955	5.9041			-go +	_ 01 _	+
Site Name: Job Number: Project Manager\Co	5812 730 (Hollis 18230		amp										ento	, CA	4 95		Ph	: 9	16.5	65.	741	12/F		910	6.56	65.7412			aroun	d	Section)
Samplers: Recorder (Signature	Required):	May	in	N	Matr	ix					ners	0	1	5											clean-up				7 1	to .	_	
Field Sample Identification No.	Date	Time	Lab Sample No.	Soil	Water	Other	_	$\overline{}$	0	90	Other	TOF	Ta	78/											Silica gel	Hold		F	Remarks			
DW-1	4-26-11	1105			X		1			X		X	2	1						-	-	+	-									
																						-	1									
							+		+	+	+	F										+		+				_				
							1	1	+		+	F					GE GOO	TI DO	0	N DI	TIO	N	+		AF	PPR	OPRIATE					
											\pm						DEC	HL	DR	INA	TEC	N	LA AS			P	RESERVED IN	LAB	3			
									+													I	I									
Relinquished by: (Signatu	156	5/	Date 4-26-1	1		7	Time	10	7	/	5	Re	cetv	ed b	y: (S	ign	ature	()		-		_	ノフ		Date	1	26/11	, [Time	71	5	-
Relinquished by: (Signati	ure)	7	Date 4/26/11			7	Time	1	20	0		Re	ceiv	ed by	y: (S	igna	ature	;)							Dale	9	2911	7	Time			
Relinquished by: (Signate	95/05/6		Date			1	Time	9				Re	ceiv	ed by	Lai	b: (S	Sign	atur	6)	1	8				Date	4	/26/11	7	Time 170	0		
Sent to Laboratory (Laboratory Commen		ME	Compbell									Me	_	Han				$\overline{}$	Pri	vate	-		ouri er (C		Nar	_	Fed Ex	A	Airborne		UPS	
	0	White Copy	- Original		Yel	low (Сор	by -	Lab	ora	tory	J				Pin	k C	ору	- F	iel	d				C	00	C Number:	01	052	46		

McCampbell Analytical, Inc.

1534 Willow Pass Rd

CHAIN-OF-CUSTODY RECORD

Page 1 of 1

Pittsburg (925) 25	g, CA 94565-1701 2-9262				•	WorkOr	der: 11	04728	(ClientCo	ode: TV	WRF				
		WaterTrax	WriteOn	☐ EDF		Excel	□F	ax	✓ Email		HardC	Сору	Third	Party	□J-f	lag
Report to: Peter Cusac	V	Email: ;	oicusack@tra	adwellrollo.com		Bil	I to:	nts Payab	ام			Requ	uested '	TAT:	1	day
Treadwell & 555 Montgon	Rollo nery St., Suite 1300 co, CA 94111	cc: PO:		5812 Hollis St			Tready 555 Mo	vell & Rol ontgomery ancisco, (lo ⁄ St., Su		ı		e Recei		04/26/ 04/26/	
								Red	uested	Tests (See leg	end be	elow)			
Lab ID	Client ID		Matrix	Collection Date	Hold	1	2 3		uested 5	Tests (See leg 7	end be	elow)	10	11	12
Lab ID 1104728-001	Client ID DW-1		Matrix Water	Collection Date 4/26/2011 11:05	Hold	1 B	2 3		•	,	See leg 7		·	10	11	12

Test Legend:

1	8260B_W	2 G-MBTEX_W	3	4	5
6		7	8	9	10
11		12			
The foll	owing SampID: 001A con	tains testgroup.			Prepared by: Ana Venegas

The following SampID: 001A contains testgroup.

Comments: SEND HARD COPY 24hr rush

> NOTE: Soil samples are discarded 60 days after results are reported unless other arrangements are made (Water samples are 30 days). Hazardous samples will be returned to client or disposed of at client expense.

Sample Receipt Checklist

Client Name:	Treadwell & Ro	lo			Date	and Time Received:	4/26/2011	5:42:33 PM
Project Name:	#730482302; 58	12 Hollis St			Chec	klist completed and re	eviewed by:	Ana Venegas
WorkOrder N°:	1104728	Matrix <u>Water</u>			Carrie	er: Rob Pringle (M	Al Courier)	
		<u>Chair</u>	of Cu	stody (C	OC) Inform	ation		
Chain of custody	present?		Yes	V	No 🗆			
Chain of custody	signed when relinqu	uished and received?	Yes	V	No \square			
Chain of custody	agrees with sample	labels?	Yes	✓	No 🗌			
Sample IDs noted	by Client on COC?		Yes	V	No \square			
Date and Time of	collection noted by C	Client on COC?	Yes	✓	No \square			
Sampler's name r	noted on COC?		Yes	V	No 🗆			
		<u>s</u>	ample	Receipt	Informatio	<u>n</u>		
Custody seals int	tact on shipping cont	ainer/cooler?	Yes		No 🗆		NA 🔽	
Shipping containe	er/cooler in good con	dition?	Yes	V	No 🗆			
Samples in prope	er containers/bottles	?	Yes	✓	No 🗆			
Sample container	rs intact?		Yes	✓	No \square			
Sufficient sample	volume for indicated	d test?	Yes	✓	No 🗌			
		Sample Prese	rvatio	n and Ho	old Time (HT	[] Information		
All samples recei	ved within holding tir	me?	Yes	✓	No 🗌			
Container/Temp B	Blank temperature		Coole	er Temp:	6.2°C		NA \square	
Water - VOA vial	ls have zero headsp	ace / no bubbles?	Yes	✓	No 🗆	No VOA vials submi	itted	
Sample labels ch	necked for correct pr	eservation?	Yes	✓	No 🗌			
Metal - pH accep	table upon receipt (p	H<2)?	Yes		No 🗆		NA 🗹	
Samples Receive	ed on Ice?		Yes	V	No 🗆			
		(Ice Typ	e: WE	TICE)			
* NOTE: If the "N	No" box is checked,	see comments below.						
=====					====	======		======
Client contacted:		Date contact	ted:			Contacted	by:	
Comments:								

Treadwell & Rollo
Client Project ID: #730482302; 5812
Hollis St
Date Sampled: 04/26/11
Date Received: 04/26/11
Client Contact: Peter Cusack
Date Extracted: 04/27/11
Client P.O.:
Date Analyzed: 04/27/11

Volatile Organics by P&T and GC/MS (Basic Target List)*

Extraction Method: SW5030B Analytical Method: SW8260B Work Order: 1104728

Extraction Method: 5 11 3030B		7 thary t	icai wicin	эч. 5 11 0200В	Work Order: 1104	720							
Lab ID		1104728-001B											
Client ID				DW-1									
Matrix				Water									
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit						
Acetone	12	1.0	10	tert-Amyl methyl ether (TAME)	ND	1.0	0.5						
Benzene	ND	1.0	0.5	Bromobenzene	ND	1.0	0.5						
Bromochloromethane	ND	1.0	0.5	Bromodichloromethane	ND	1.0	0.5						
Bromoform	ND	1.0	0.5	Bromomethane	ND	1.0	0.5						
2-Butanone (MEK)	ND	1.0	2.0	t-Butyl alcohol (TBA)	ND	1.0	2.0						
n-Butyl benzene	ND	1.0	0.5	sec-Butyl benzene	ND	1.0	0.5						
tert-Butyl benzene	ND	1.0	0.5	Carbon Disulfide	ND	1.0	0.5						
Carbon Tetrachloride	ND	1.0	0.5	Chlorobenzene	ND	1.0	0.5						
Chloroethane	ND	1.0	0.5	Chloroform	ND	1.0	0.5						
Chloromethane	ND	1.0	0.5	2-Chlorotoluene	ND	1.0	0.5						
4-Chlorotoluene	ND	1.0	0.5	Dibromochloromethane	ND	1.0	0.5						
1,2-Dibromo-3-chloropropane	ND	1.0	0.2	1,2-Dibromoethane (EDB)	ND	1.0	0.5						
Dibromomethane	ND	1.0	0.5	1.2-Dichlorobenzene	ND	1.0	0.5						
1.3-Dichlorobenzene	ND	1.0	0.5	1.4-Dichlorobenzene	ND	1.0	0.5						
Dichlorodifluoromethane	ND	1.0	0.5	1,1-Dichloroethane	ND	1.0	0.5						
1,2-Dichloroethane (1,2-DCA)	ND	1.0	0.5	1,1-Dichloroethene	ND	1.0	0.5						
cis-1,2-Dichloroethene	ND	1.0	0.5	trans-1,2-Dichloroethene	ND	1.0	0.5						
1,2-Dichloropropane	ND	1.0	0.5	1,3-Dichloropropane	ND	1.0	0.5						
2,2-Dichloropropane	ND	1.0	0.5	1,1-Dichloropropene	ND	1.0	0.5						
cis-1,3-Dichloropropene	ND	1.0	0.5	trans-1,3-Dichloropropene	ND	1.0	0.5						
Diisopropyl ether (DIPE)	10	1.0	0.5	Ethylbenzene	ND	1.0	0.5						
Ethyl tert-butyl ether (ETBE)	ND	1.0	0.5	Freon 113	ND	1.0	10						
Hexachlorobutadiene	ND	1.0	0.5	Hexachloroethane	ND	1.0	0.5						
2-Hexanone	ND	1.0	0.5	Isopropylbenzene	ND	1.0	0.5						
4-Isopropyl toluene	ND	1.0	0.5	Methyl-t-butyl ether (MTBE)	ND	1.0	0.5						
Methylene chloride	ND	1.0	0.5	4-Methyl-2-pentanone (MIBK)	ND	1.0	0.5						
Naphthalene	ND	1.0	0.5	n-Propyl benzene	ND	1.0	0.5						
Styrene	ND	1.0	0.5	1,1,1,2-Tetrachloroethane	ND	1.0	0.5						
1,1,2,2-Tetrachloroethane	ND	1.0	0.5	Tetrachloroethene	ND	1.0	0.5						
Toluene	ND	1.0	0.5	1,2,3-Trichlorobenzene	ND	1.0	0.5						
1.2.4-Trichlorobenzene	ND	1.0	0.5	1.1.1-Trichloroethane	ND	1.0	0.5						
1,1,2-Trichloroethane	ND	1.0	0.5	Trichloroethene	0.57	1.0	0.5						
Trichlorofluoromethane	ND	1.0	0.5	1,2,3-Trichloropropane	ND	1.0	0.5						
1.2.4-Trimethylbenzene	ND	1.0	0.5	1,3,5-Trimethylbenzene	ND	1.0	0.5						
Vinyl Chloride	ND	1.0	0.5	Xylenes, Total	ND	1.0	0.5						
		Surr		ecoveries (%)			***						
%SS1:	9			%SS2:	10	15							
%SS3:	8	/0332.	10	13									
/0000.	1 0	U		1									

Comments

ND means not detected above the reporting limit/method detection limit; N/A means analyte not applicable to this analysis; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

surrogate diluted out of range or coelutes with another peak; &) low surrogate due to matrix interference.

^{*} water and vapor samples are reported in $\mu g/L$, soil/sludge/solid samples in mg/kg, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts are reported in mg/L, wipe samples in $\mu g/W$ in μg

McCampbell Analytical, Inc. "When Ouality Counts"

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo	Client Project ID: #730482302; 5812 Hollis St	Date Sampled: 04/26/11
555 Montgomery St., Suite 1300	Hollis St	Date Received: 04/26/11
	Client Contact: Peter Cusack	Date Extracted: 04/27/11
San Francisco, CA 94111	Client P.O.:	Date Analyzed 04/27/11

Gasoline Range (C6-C12) Volatile Hydrocarbons as Gasoline*

Extraction method	SW5030B	Analytic	al methods SW8015Bm	Wo	rk Order:	1104728	
Lab ID	Client ID	Matrix	TPH(g)	DF	% SS	Comments	
001A	DW-1	W	ND	1	106		
				<u> </u>			
	porting Limit for DF =1; means not detected at or	W	50		μg/L		
	ove the reporting limit	S	NA		NA		

^{*} water and vapor samples are reported in ug/L, soil/sludge/solid samples in mg/kg, wipe samples in μ g/wipe, product/oil/non-aqueous liquid samples and all TCLP & SPLP extracts in mg/L.

Angela Rydelius, Lab Manager

[#] cluttered chromatogram; sample peak coelutes w/surrogate peak; low surrogate recovery due to matrix interference. %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

McCampbell Analytical, Inc. "When Quality Counts"

1534 Willow Pass Road, Pittsburg, CA 94565-1701 Web: www.mccampbell.com E-mail: main@mccampbell.com Telephone: 877-252-9262 Fax: 925-252-9269

Treadwell & Rollo		ID: #730482302; 5812		Date Sampled:	04/26/1	11		
555 Montgome	ery St., Suite 1300	Hollis St			Date Received:	04/26/1	11	
333 Montgonie	ery St., Suite 1500	Client Contac	t: Peter Cusack		Date Extracted:	04/26/1	11	
San Francisco,	CA 94111	Client P.O.:		04/27/11				
Extraction method:	SW3510C		ctable Petroleum Hydroc	Work Order: 1104728				
Lab ID	Client ID	Matrix	TDH Diasal TDH Motor O				% SS	Comments
1104728-001A	DW-1	W	ND		ND	1	92	

* water samples are reported in µg/L, wipe sa	imples in µg/wipe, soil/solid/sludg	e samples in mg/kg, produc	t/oil/non-aqueous liquid s	samples in mg/L, a	and all
DISTLC / STLC / SPLP / TCLP extracts are	reported in ug/L.				

50

NA

W

S

cluttered chromatogram resulting in coeluted surrogate and sample peaks, or; surrogate peak is on elevated baseline, or; surrogate has been diminished by dilution of original extract; %SS = Percent Recovery of Surrogate Standard; DF = Dilution Factor

+The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation:

DHS ELAP Certification 1644

Reporting Limit for DF =1;

ND means not detected at or

above the reporting limit

Angela Rydelius, Lab Manager

250

NA

μg/L

mg/Kg

QC SUMMARY REPORT FOR SW8260B

W.O. Sample Matrix: Water QC Matrix: Water BatchID: 57862 WorkOrder 1104728

EPA Method SW8260B	Extra	ction SW	5030B					S	piked San	nple ID	: 1104692-0	007B
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%))
, and y	μg/L	μg/L	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
tert-Amyl methyl ether (TAME)	ND	10	89.2	87.9	1.47	93.8	90	4.08	70 - 130	30	70 - 130	30
Benzene	ND	10	98.3	97.3	1.05	112	108	4.39	70 - 130	30	70 - 130	30
t-Butyl alcohol (TBA)	ND	50	95.5	90.9	4.93	94.5	93.8	0.833	70 - 130	30	70 - 130	30
Chlorobenzene	ND	10	110	106	3.43	128	118	8.46	70 - 130	30	70 - 130	30
1,2-Dibromoethane (EDB)	ND	10	111	109	2.29	120	110	8.91	70 - 130	30	70 - 130	30
1,2-Dichloroethane (1,2-DCA)	ND	10	102	100	1.59	111	106	4.44	70 - 130	30	70 - 130	30
1,1-Dichloroethene	ND	10	106	103	2.36	123	118	4.08	70 - 130	30	70 - 130	30
Diisopropyl ether (DIPE)	ND	10	105	105	0	118	114	3.62	70 - 130	30	70 - 130	30
Ethyl tert-butyl ether (ETBE)	ND	10	98.3	98.1	0.196	107	103	3.35	70 - 130	30	70 - 130	30
Methyl-t-butyl ether (MTBE)	0.58	10	102	101	0.941	106	105	1.42	70 - 130	30	70 - 130	30
Toluene	ND	10	102	99.4	2.68	116	109	6.54	70 - 130	30	70 - 130	30
Trichloroethene	ND	10	95.6	93.8	1.92	111	105	5.34	70 - 130	30	70 - 130	30
%SS1:	92	25	98	97	1.08	96	97	1.39	70 - 130	30	70 - 130	30
%SS2:	100	25	100	100	0	100	100	0	70 - 130	30	70 - 130	30
%SS3:	87	2.5	101	100	0.741	97	98	1.50	70 - 130	30	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57862 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104728-001B	04/26/11 11:05 AM	04/27/11	04/27/11 8:19 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

surrogate diluted out of range or coelutes with another peak; &) low surrogate due to matrix interference.

Laboratory extraction solvents such as methylene chloride and acetone may occasionally appear in the method blank at low levels.

QA/QC Officer

QC SUMMARY REPORT FOR SW8021B/8015Bm

W.O. Sample Matrix: Water QC Matrix: Water BatchID: 57895 WorkOrder 1104728

EPA Method SW8015Bm	Extra	ction SW	5030B					S	Spiked San	nple ID	: 1104704-0	05A
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
7 mary to	μg/L	μg/L	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH(btex [£]	ND	60	126	110	13.5	122	125	2.82	70 - 130	20	70 - 130	20
MTBE	ND	10	97.2	98	0.731	94.2	99.1	5.06	70 - 130	20	70 - 130	20
Benzene	ND	10	96.8	90.9	6.30	93.9	94.7	0.884	70 - 130	20	70 - 130	20
Toluene	ND	10	95.5	89.2	6.75	92.9	93.4	0.451	70 - 130	20	70 - 130	20
Ethylbenzene	ND	10	97	90.2	7.30	94.3	95	0.676	70 - 130	20	70 - 130	20
Xylenes	ND	30	97.1	91.5	5.94	94.5	95.2	0.665	70 - 130	20	70 - 130	20
%SS:	98	10	99	89	10.3	98	98	0	70 - 130	20	70 - 130	20

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57895 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed
1104728-001A	04/26/11 11:05 AM	1 04/27/11	04/27/11 10:59 AM				

MS = Matrix Spike; MSD = Matrix Spike Duplicate; LCS = Laboratory Control Sample; LCSD = Laboratory Control Sample Duplicate; RPD = Relative Percent Deviation.

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

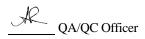
MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

£ TPH(btex) = sum of BTEX areas from the FID.

cluttered chromatogram; sample peak coelutes with surrogate peak.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = matrix interference and/or analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content, or inconsistency in sample containers.



QC SUMMARY REPORT FOR SW8015B

W.O. Sample Matrix: Water QC Matrix: Water BatchID: 57832 WorkOrder 1104728

EPA Method SW8015B					5	Spiked San	nple ID:	N/A				
Analyte	Sample	Spiked	MS	MSD	MS-MSD	LCS	LCSD	LCS-LCSD	Acce	eptance	Criteria (%)	
7 may to	μg/L	μg/L	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	MS / MSD	RPD	LCS/LCSD	RPD
TPH-Diesel (C10-C23)	N/A	1000	N/A	N/A	N/A	105	105	0	N/A	N/A	70 - 130	30
%SS:	N/A	625	N/A	N/A	N/A	95	95	0	N/A	N/A	70 - 130	30

All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions: NONE

BATCH 57832 SUMMARY

Lab ID	Date Sampled	Date Extracted	Date Analyzed	Lab ID	Date Sampled	Date Extracted	Date Analyzed	
1104728-001A	04/26/11 11:05 AM	1 04/26/11	04/27/11 12:21 AM					

 $MS = Matrix \; Spike; \\ MSD = Matrix \; Spike \; Duplicate; \\ LCS = Laboratory \; Control \; Sample; \\ LCSD = Laboratory \; Control \; Sample \; Duplicate; \\ RPD = Relative \; Percent \; Deviation.$

% Recovery = 100 * (MS-Sample) / (Amount Spiked); RPD = 100 * (MS - MSD) / ((MS + MSD) / 2).

MS / MSD spike recoveries and / or %RPD may fall outside of laboratory acceptance criteria due to one or more of the following reasons: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) the spiked sample's matrix interferes with the spike recovery.

N/A = not enough sample to perform matrix spike and matrix spike duplicate.

NR = analyte concentration in sample exceeds spike amount for soil matrix or exceeds 2x spike amount for water matrix or sample diluted due to high matrix or analyte content.

QA/QC Officer



Appendix B

March 2010 NPDES Self-Monitoring Report

NPDES SELF-MONITORING REPORT TREATMENT SYSTEM STARTUP FIRST QUARTER 2011 GENERAL PERMIT CAG912002 SFRWQCB ORDER R2-2006-0075 Construction Dewatering Treatment System Greenway Phase II 5812 Hollis Street Emeryville, California

> Geotracker Site ID T0600101109 CIWQS Place ID 753920

> > March 2010

Prepared on Behalf of

Pacific States Environmental Contractors, Inc. 11555 Dublin Boulevard Dublin, California 94568

Prepared for

EmeryStation Triangle 2, LLC Greenway Phase II 1120 Nye Street, Suite 400 San Rafael, CA 94901

Prepared by



3815 Brighton Avenue, Oakland, California 94602 broat@earthlink.net, 510.919.4358

Rev. 0 - March 30, 2011



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SIGNATURE PAGE

All engineering information, conclusions, and recommendations contained in this report have been prepared by a California Professional Engineer.

Robert E. Roat, P.E.

California Professional Engineer (53593)

3/31/4/



1.0 INTRODUCTION

This NPDES Start-up and First Quarter 2011 Self-Monitoring Report (SMR) has been prepared by Brighton Environmental Consulting to present the startup results and first quarter 2011 results of the National Pollutant Discharge Elimination System (NPDES) Self-Monitoring Program for the temporary construction dewatering and treatment system ("the Treatment System") at 5812 Hollis Street in Emeryville, California. Treated groundwater from dewatering activities is being discharged to the Caltrans-operated storm drain system under the *GENERAL WASTE DISCHARGE REQUIREMENTS FOR Discharge or Reuse of Extracted and Treated Groundwater Resulting from the Cleanup of Groundwater Polluted by Fuel Leaks and Other Related Wastes at Service Stations and Similar Sites (San Francisco Bay Regional Water Quality Control Board Order NO. R2-2006-0075, National Pollutant Discharge Elimination System (NPDES) PERMIT NO. CAG912002*, herein referred to as the NPDES Permit. The Owner has submitted a Notice of Intent for coverage under the NPDES Permit (Greenway 2010) and has received a letter of authorization from the San Francisco Bay Regional Water Quality Control Board (Regional Board 2010).

This report summarized activities, sampling and laboratory analyses results associated with the startup and operation of the treatment system in March 2011.

1.1 COMPLIANCE SUMMARY

1.1.1 Effluent Limitations

The NPDES Permit requires that the discharge of the effluent maintain compliance with the effluent limitations at a discharge point after full treatment but before it joins or is diluted by any other waste stream, body of water, or substance. The effluent limits for specific compounds are defined in the permit. The effluent limits for the organic compounds listed in the permit and for pH (NPDES Permit Section IV.A) were not exceeded during start up. Average daily flow rate



did not exceed the daily flow rate of 100 gpm (144,000 gallons per day) permitted by the letter of authorization.

1.1.2 Trigger Levels

Triggers are not effluent limitations; instead, they are levels at which additional investigation is warranted to determine whether a numeric limit for a particular constituent is necessary. If any constituent in the effluent of a discharge exceeds the trigger levels, then additional sampling, analyses, evaluation and/or treatment for those constituents are required by the NPDES Permit.

The concentration-based triggers for volatile organic compounds, semi-volatile organic compounds and inorganic compounds tested were not exceeded in effluent samples during startup (NPDES Permit Section 6.C.6), with the following exceptions:

- total copper at 6.2 μg/L, above the trigger level of 3.1 μg/L (February 28,2011).
- MDLs for cadmium exceeded the trigger level of 0.07 μg/L
- TEPH as motor oil reporting limit exceeds the trigger level of 50 µg/L. J-flagged concentration in one sample was at the trigger level, although the samples were also flagged indicating concentrations of TEPH in the laboratory blank.

In addition, di-isopropyl ether (DIPE) was reported in influent samples above reporting limits. Monthly VOC sampling will be conducted to monitor for DIPE in influent, midpoint and effluent samples.

1.2 SITE LOCATION

The project is located 5812 Hollis Street in Emeryville, California (Figure 1-1).

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1.3 Purpose of the Groundwater Discharge

ES Triangle 2 LLC is developing a multistory office building. As part of the construction, the

site will be excavated to allow installation of subsurface facilities. A dewatering system will be

installed to control groundwater flow into the excavation. The dewatering system will be

operated for approximately 12 months during the excavation and installation of subsurface

facilities.

1.4 CHEMICALS OF CONCERN

Previous investigations at the site have revealed the presence of petroleum hydrocarbons and

chlorinated solvents in groundwater. The site is currently being investigated and remediated by

the Owner (Geotracker Global ID T0600101109).

Because concentrations of petroleum hydrocarbons and chlorinated solvents may be present in

pumped groundwater, PSEC plans to collect and treat dewatering system effluent prior to

discharge to the storm drain system.

As summarized below, results of startup sampling have revealed the presence of chlorinated

solvents and fuel additives in influent samples. The concentrations in influent samples were

below effluent limitations.

1.5 RESPONSIBLE PARTIES

The property and facility Owner is ES Triangle 2 LLC. The project General Contractor is DPR

Construction, Inc. of Redwood City, California. PSEC will provide, maintain and inspect

treatment system equipment. Viking Drillers, Inc. of West Sacramento, California will install

and maintain dewatering wells. Brighton and/or PSEC will perform daily inspections and flow

monitoring and compliance sampling and analyses. The Civil Engineer providing technical

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oversight and compliance monitoring of the Treatment System operation is Brighton Environmental Consulting of Oakland, California.

1.6 NOTIFICATION OF NPDES COVERAGE

The Owner submitted to the Regional Board a Notice of Intent for coverage under the NPDES Permit in June 2010 (Greenway 2010). After review the NOI, The Regional Board issued a *Notice of General Permit Coverage*, under the requirements of the NPDES Permit (NGPC letter, Regional Board, July 12, 2010). The NGPC letter allowed discharge under the NPDES permit with the additional conditions as summarized below:

- 1. Use CIWQS Place identification number and the facility address listed in the table for any future communications.
- 2. Notify Lou Gonzales within 24 hours at the e-mail address below, of the document titles uploaded on Geotracker.
- 3. Submit the Self-Monitoring Reports no later than 45 days following the last day of the quarter, and the Annual Reports by February 15th of each year. If there is no discharge during the reporting period, please state this information in the reports. Late reports and effluent violations are subject to mandatory minimum penalties pursuant to California Water Code 13385 (h)&(i).
- 4. Operate the dewatering treatment system 24/7 during its approximately seventeen-month duration. The dewatering treatment system will not be allowed to discharge to storm drain if the treatment system is not operable.
- 5. Make available the back-up treatment system components on a 24-hour turnaround time.
- 6. Conduct daily, weekly and monthly inspections of the treatment system by designated

1.7 PURPOSE OF THE REPORT

This SMR presents information gathered during the start-up operation of the Treatment System between March 3, 2011 and March 11, 2011, as well as flow information through the end of March 2011.

The Site location and the discharge point to Temescal Creek and the San Francisco Bay are shown on Figure 1-1. A process flow diagram of the Treatment System showing the Self-Monitoring Program sampling points is presented on Figure 1-2.

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1.8 DESCRIPTION OF TREATMENT SYSTEM

Groundwater from 20 wells is pumped through a collection system to one 20,000-gallon baffled

settling tank (Settling Tank 1) to remove suspended solids. Water then flows by gravity from an

outlet on the upper part of the settling tank to a 20,000-gallon batch tank. Using a submersible

pump, water from the batch tank is transferred through a two-pod sand filter system to remove

additional suspended solids. Water is then pumped through two 1,000-pound activated carbon

adsorption units arranged in series. After carbon treatment, the water is discharged through a

totalizing flow meter to the storm drain. Sand filter cleaning occurs as needed based on pressure

differential and flow rate losses. Carbon changeout will occur based on analytical results for VOCs

and TPH from a midpoint sample.

The treatment system submersible pump was capable of transferring at least 100 gallons per

minute at 50 feet of pressure. The process pump flow rate is adjusted to maximize dewatering,

but is regulated so as not to exceed a flow rate of 100 gallons per minute.

System controls include level switches in the SRT. A high level switch in the batch tank

activates to turn on the system transfer pump. A low level switch in the batch tank activates to

turn off the system transfer pump. A high-high level switch activates to shut off power to the

well pumps if the Treatment System pumps are not operating or not pumping at a sufficient rate

to support dewatering flow. A separate low level switch in the batch tanks turns the dewatering

well pumps on when sufficient freeboard is available.

After discharge from the activated carbon adsorption vessel, treated water flows to a storm drain

inlet at the northwest corner of Hollis Street and Powell Street. The discharge path for the storm

drain is shown in Figure 1-1.

1.8.1 Deviations from the NOI

The system was installed and operated as described in the NOI.

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2.0 TREATMENT SYSTEM OPERATION AND MONITORING

2.1 SUMMARY OF SYSTEM OPERATION

The dewatering wells and treatment system were constructed in February 2011. On February 28, 2011, the dewatering system began pumping to allow testing of the treatment system without discharge. Water was pumped to the settling tank and batch tank and the treatment system began pumping, with water recycled to the settling tank. During initial pumping, influent, effluent and midpoint samples were collected and analyzed for the constituents shown in Appendix A. Water was then stored and recycled through the treatment system until startup. Treatment system discharge to the storm drain began after receipt of positive analytical results on March 3, 2011.

2.2 STARTUP INSPECTIONS

The treatment system was inspected several times by the Engineer during startup: 1) Prior to initial testing on February 28, 2011, 2) during start of discharge on March 3, 2011, and during the second startup sampling event on March 7, 2011. The system was inspected for conformance to plans and to best practices, for structural integrity, and for potential failure modes. Several issues were identified and corrected as a result of these inspections, including

- 1. Installation of the high-high level well shutoff
- 2. The need for an influent sample port
- 3. The need to label sample ports
- 4. A small leak in a fitting at the submersible pump coupling on top of the batch tank, dripping back into the batch tank

Items 1 through 3 were addressed before system start up. The fourth issue was addressed by March 11, 2011.



2.3 SAMPLING AND ANALYSIS METHODS

2.3.1 List of Approved Analyses

All laboratory analyses were performed by commercial analytical laboratories certified by the State of California Department of Health Services to perform the particular analyses. Summaries of analyses, quality control measures, and laboratories used are included in Appendix A.

2.3.2 Sample Locations

Pursuant to the NPDES Self-Monitoring Program, the NGPC letter and the General Permit requirements, water samples can be collected for laboratory analysis from Treatment System sample ports (Figure 1-2) and from a receiving water location at the storm drain system outlet (Figure 1-1).

Sample locations for the Self-Monitoring Program for system startup included compliance sampling stations INF-001 (influent), EFF-001 (effluent), and operational sampling station MID-001. Samples were collected by trained sample collection staff of PSEC.

2.3.3 Sampling Methods

Influent samples (INF-001) were collected directly from the inlet pipe at the settling tank. Samples from MID-001 were collected from a sample tap at the effluent of the first carbon adsorption unit. EFF-001 samples were collected from a sample tap at the effluent of the second carbon adsorption unit. Samples were collected directly into laboratory bottles with the required preservative, if needed. Prior to sample collection the pH, electrical conductivity and temperature were analyzed in the field with an YSI 63 pH/conductivity meter, and the turbidity was measured with a Hach 2100P turbidity meter. Samples were not collected until three consecutive samples were within five percent of each other for each field analysis.



After filling and labeling, the sample containers were placed in chilled, insulated coolers for transport to the laboratory for analysis. Chain-of-custody forms were completed for the samples. These forms accompanied the samples until receipt by the California Department of Health Services-certified laboratory listed in Appendix A. Samples from the first day sampling were analyzed on a same day turn around. The system was allowed to begin discharge after the Engineer reviewed the analytical results and approved the effectiveness of treatment. Samples from the five day sampling were also analyzed on a same day turn around, due to lab shutdowns during the Thanksgiving holiday.

2.3.4 Laboratory Analysis

During startup, samples were collected and analyzed for parameters shown in Table 2-1. Laboratory quality assurance/quality control (QA/QC) data and reporting limits were reviewed for each laboratory report received. Results of the review are summarized in Appendix A.



3.0 RESULTS

3.1 FIELD MEASUREMENT RESULTS

3.1.1 Standard Observations

Standard observations of the effluent were made at EFF-001 on March 3, 2011 and March 7, 2011. A summary of the observations is included in Appendix B.

3.1.2 Flow Rate and Volume

Average flow rate for the startup period (March 3, 2011 through March 11, 2011) was calculated at approximately 5.4 gallons per minute (gpm), or approximately 7,829 gallons per day (gpd). The total volume of groundwater pumped, treated and discharged during startup was approximately 0.065 million gallons. Average flow rate for the First Quarter of 2011 (March 3, 2011 through March 28, 2011) was calculated at approximately 8.5 gallons per minute (gpm), or approximately 12,240 gallons per day (gpd). The total volume of groundwater pumped, treated and discharged during startup was approximately 0.31 million gallons. Flow rates and daily volumes are summarized in Table 3-1.

The average daily flow rate was below the maximum permitted rate of 144,000 gpd (100 gpm).

3.2 LABORATORY ANALYTICAL RESULTS

The Fuel General NPDES Permit contains effluent limitations for 21 organic pollutants, pH and toxicity (CAG 912002.IV.A.1). The Permit also contains trigger concentrations for numerous organic compounds and metals.

Results of self-monitoring analyses for system startup are summarized in Tables 3-2 through 3-6 and discussed below. Quality assurance results for the analyses are summarized in Appendix A.

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3.2.1 Turbidity and pH

Water samples from sampling stations INF-001 and EFF-001 were collected and analyzed in the

laboratory for turbidity by Standard Method 2130B and pH by EPA Method 9040C. Results are

included in Table 3-2.

During startup, the pH of the effluent water was within the effluent limit range specified in the

NPDES permit. Turbidity in the effluent samples exceeded the trigger level for field samples but

not laboratory analyses.

3.2.2 Volatile Organic Compounds with Effluent Limitations

Water samples were analyzed for volatile organic compounds (VOCs) by EPA Method 8260B.

Laboratory analyses of effluent samples, midpoint samples and influent samples collected during

startup revealed no VOCs with concentrations above effluent limitations or laboratory reporting

limits.

As required by the NPDES permit, the laboratory also reported "J-flagged" results, indicating

that the laboratory observed concentrations of the compound above method detection limits

(MDLs) but below laboratory reporting limits. No VOCs with effluent limits were reported

above MDLs for the EFF-001 or MID-001 samples. Toluene, trichloroethene and vinyl chloride

were reported in INF-001 samples at concentrations below reporting limits but above method

detection limits. Analytical results are shown in Table 3-3 for organic compounds with permit-

specified numerical effluent limitations.

3.2.3 Total Petroleum Hydrocarbons

Water samples from sampling stations INF-001 and EFF-001 were collected for analysis for

TPH as gasoline and TEPH as diesel fuel and motor oil. Water samples were also collected from

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MID-001 for TPH as gasoline. TPH as gasoline was analyzed by EPA 8260B and TPH as diesel fuel and motor oil were analyzed by EPA Method 8015M.

Concentrations of TPH as gasoline and TEPH as diesel fuel in effluent samples were below effluent limitations, laboratory reporting limits and MDLs during startup sampling. Concentrations of TPH as gasoline in midpoint samples were below effluent limitations, laboratory reporting limits and MDLs during startup sampling. Laboratory analyses for one influent sample revealed the presence of TEPH as diesel fuel at 24 μ g/L, below the reporting limit but above the method detection limit, although the laboratory also reported TEPH as diesel in the laboratory blank. Results of TPH as gasoline and TEPH as diesel fuel are summarized in Table 3-3.

For EFF-001 samples, concentrations of TEPH as motor oil were below or at trigger concentrations, but below reporting limits during startup sampling. For INF-001 samples, concentrations of TEPH as motor oil ranged from ND to 51 ug/L (J-flagged) during startup sampling. Results of TEPH as motor oil are summarized in Table 3-4.

3.2.4 Volatile Organic Compounds and Volatile Fuel Additives with Trigger Levels

Laboratory analyses for VOCs with numerical trigger levels revealed no VOCs with trigger levels in effluent (EFF-001) startup samples at concentrations above laboratory reporting limits. EFF-001 samples conteinaed carbon disulfide and 1,2,4-trimethyl benzene at concentrations above method detection limits but below laboratory reporting limits. None of the fuel additives were reported above MDLs in the EFF-001 sample. The laboratory did report di-isopropyl ether (DIPE), a fuel additive above reporting limits in both INF-001 samples. Laboratory analytical results for VOC compounds with reported concentrations above MDLs and for volatile fuel additives are shown in Table 3-4. Due to the presence of DIPE, additional analyses for fuel oxygenates and VOCs will be conducted in Second Quarter 2011.

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3.2.5 Ethylene Dibromide

Water samples from sampling stations INF-001 and EFF-001 were collected for analysis on

March 7, 2011 and analyzed for ethylene dibromide (EDB) by EPA Method 504.1. Analytical

results for EDB and other VOCs are summarized in Table 3-3. EDB was not reported above

MDLs in either influent or effluent samples.

3.2.6 Semi Volatile Organic Compounds

Water samples from INF-001 and EFF-001 were collected on February 28, 2011 and analyzed

for semi-volatile organic compounds (sVOCs) by EPA Method 8270C. Concentrations of

sVOCs in effluent samples were below laboratory reporting limits and MDLs for all analytes.

Bis (2-ethyl hexyl phthalate was reported at 3.7 μg/L, below the trigger level of 5 μg/L, in the

INF-001 sample. Results of laboratory analyses for sVOCs are summarized in Table 3-4.

3.2.7 Polynuclear Aromatic Hydrocarbons

Water samples from INF-001 and EFF-001 were collected on March 7, 2011 and analyzed for

polynuclear aromatic hydrocarbons (PAHs) by EPA Method 8310. Concentrations of PAHs in

influent and effluent samples were below laboratory reporting limits and trigger levels for all

analytes. Flourene was reported above MDLs in the INF-001 sample. Results of laboratory

analyses for PAHs are summarized in Table 3-4.

3.2.8 Inorganic Analyses

Samples were collected during both startup sampling events and analyzed for total metals and

total cyanide. Inorganic constituents analyzed by EPA Method 6020 or 7000 series tests during

startup included antimony, arsenic, beryllium, cadmium, total chromium, copper, lead, nickel,

selenium, silver, thallium, and zinc. Mercury was analyzed by EPA Method 7470. Low level

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mercury was analyzed by EPA Method 1631 for samples from March 7, 2011. Total cyanide was analyzed using standard method 4500CN-E. Results are summarized in Table 3-5.

Copper was reported in one EFF-001 sample at concentrations in excess of the trigger level. Chromium, lead and nickel were reported in at least one EFF-001 sample at concentrations above reporting limits but below trigger levels. Antimony, selenium and zinc were reported in at least one EFF-001 sample above MDLs but below reporting limits. Cadmium MDLs exceeded the trigger concentration. Mercury MDLs exceeded trigger levels, but analyses for low-level mercury revealed that concentrations of total mercury are below trigger levels. Total cyanide was not reported at concentrations above laboratory reporting limits or trigger levels. Additional analyses for cadmium and copper will be conducted during second quarter of 2011.

3.2.9 Mass Removal of Organic Compounds

Approximately 0.0008 pounds of organic compounds were removed from extracted groundwater during startup. Mass removal calculations for VOCs are summarized in Table 3-6. The calculations are based on the average concentrations and the total volume of groundwater extracted during startup. Only compounds with concentrations above method detection limits are included in the calculations.



mercury was analyzed by EPA Method 1631 for samples from March 7, 2011. Total cyanide was analyzed using standard method 4500CN-E. Results are summarized in Table 3-5.

Copper was reported in one EFF-001 sample at concentrations in excess of the trigger level. Chromium, lead and nickel were reported in at least one EFF-001 sample at concentrations above reporting limits but below trigger levels. Antimony, selenium and zinc were reported in at least one EFF-001 sample above MDLs but below reporting limits. Cadmium MDLs exceeded the trigger concentration. Mercury MDLs exceeded trigger levels, but analyses for low-level mercury revealed that concentrations of total mercury are below trigger levels. Total cyanide was not reported at concentrations above laboratory reporting limits or trigger levels. Additional analyses for cadmium and copper will be conducted during second quarter of 2011.

3.2.9 Mass Removal of Organic Compounds

Approximately 0.0008 pounds of organic compounds were removed from extracted groundwater during startup. Approximately 0.0039 pounds of organic compounds were removed from extracted groundwater during First Quarter 2011. Mass removal calculations for VOCs are summarized in Table 3-6. The calculations are based on the average concentrations and the total volume of groundwater extracted during startup. Only compounds with concentrations above method detection limits are included in the calculations.



5.0 REFERENCES

- Greenway 2010. Attachment B Notice of Intent Application Form to Receive Authorization to Discharge Treated Groundwater under the Requirements of the Order No. R2-2006-0075, NPDES Permit No. CAG912002 (Fuel), Greenway Phase II, 5812 Hollis Street, Emeryville, CA 94608, June 2010.
- Brighton, 2010. Attachment 2 to NOI: Engineering Certification Report, Construction Dewatering Treatment System, Greenway Phase II, 5812 Hollis Street, Emeryville, California, Geotracker Site IDs T0600101109, 60001261, June 2010.
- Regional Board (California Regional Water Quality Control Board, San Francisco Bay Region), 2006. Order No. R2-2006-0075, NPDES Permit Number CAG912002, General Waste Discharge Requirements for Discharge or Reuse of Extracted and Treated Groundwater Resulting from the Cleanup of Groundwater Polluted by Fuel Leaks and Other Related Wastes at Service Stations and Similar Sites, 2006.
- Regional Board 2010. Notice of General Permit Coverage for Discharge from 5812 Hollis Street, Emeryville, Alameda County, CA 94608, under the Requirements of Order No. R2-2006-0075, NPDES Permit No. CAG912002 (Fuels General Permit), July 12, 2010.



TABLES

							arter of ation	2nd Q	uarter of Ope	eration	3rd Q	uarter of Ope	eration
Type of Sample	Analytical Method	INF-001	MID-001	EFF-001	RSW-001U/ RSW-001D	Star	rtup	Apr-11	May-11	Jun-11	Jul-11	Aug-11	Sep-11
						Day 1	Day 5						
Bioassay 96 hour static renewal survival w/ rainbow trout	EPA-821-R-02- 012			Q/Y				EFF			EFF		
Standard	Field			М				EFF	EFF	EFF	EFF	EFF	EFF
Observations	Measurement				V								
VOC method		Y											
8020 (includes	BTEX -EPA		Y										
BTEX and MTBE) (if not running	Method 8020 MTBE - EPA Method 8260B			Y									
EPA8260B)					V								
VOC method		D/Q											
8020 (includes	BTEX -EPA		D/M			1	-				1		
BTEX only) (if not running	Method 8020			D/M									
EPA8260B)					V								
		D/Q				INF	INF	INF			INF		
TPHg (8015M)	EPA Method 8015 Mod			D/M		EFF	EFF	EFF	EFF	EFF	EFF	EFF	EFF
					V								
TPHd (8015M)	EPA Method	D/Q				INF	INF	INF	INF(T)	INF(T)	INF		
with silica gel cleanup (diesel	8015 Mod with silica gel			D/M		EFF	EFF	EFF	EFF	EFF	EFF	EFF	EFF
and motor oil)	cleanup				V	-	-				-		

						4th Q	uarter of Ope	eration	5th Q	uarter of Ope	eration
Type of Sample	Analytical Method	INF-001	MID-001	EFF-001	RSW-001U/ RSW-001D	Oct-11	Nov-11	Dec-11	Jan-12	Feb-12	Mar-12
Bioassay 96 hour static renewal survival w/ rainbow trout	EPA-821-R-02- 012			Q/Y		EFF		-1	EFF		
Standard	Field			М		EFF	EFF	EFF	EFF	EFF	EFF
Observations	Measurement				V						
VOC method		Y									
8020 (includes	BTEX -EPA		Y								
BTEX and MTBE) (if not running	Method 8020 MTBE - EPA Method 8260B			Y							
EPA8260B)					V						
VOC method		D/Q		-		1		1	1		
8020 (includes	BTEX -EPA		D/M								
BTEX only) (if not running	Method 8020			D/M							
EPA8260B)					V						
		D/Q				INF			INF		
TPHg (8015M)	EPA Method 8015 Mod			D/M		EFF	EFF	EFF	EFF	EFF	EFF
					V						
TPHd (8015M)	EPA Method	D/Q				INF			INF		
with silica gel cleanup (diesel	8015 Mod with silica gel			D/M		EFF	EFF	EFF	EFF	EFF	EFF
and motor oil)	cleanup				V						

						1st Qua Oper	arter of ation	2nd Q	uarter of Ope	eration	3rd Q	uarter of Ope	eration
Type of Sample	Analytical Method	INF-001	MID-001	EFF-001	RSW-001U/ RSW-001D	Star	tup	Apr-11	May-11	Jun-11	Jul-11	Aug-11	Sep-11
						Day 1	Day 5						
		Q					INF			INF			INF
PAH method 8310	EPA Method 8310			Q			EFF			EFF			EFF
					V								
		Y					INF						
EDB method 504	EPA Method 504			Y			EFF						
301	301				V								
		Y				INF	INF	INF	INF	INF	INF	INF	INF
VOC method	EPA Method		Y			MID	MID	MID	MID	MID	MID	MID	MID
8260b	8260B			Y		EFF	EFF	EFF	EFF	EFF	EFF	EFF	EFF
					V								
		Y					INF						
Fuel Oxygenates TAME, DIPE, ETBE, TBA, Ethanol, Methanol (with	EPA Method 8260B			Q (annual if not reported above 5 ug/l in first sample)			EFF			EFF			
VOC analyses)					V								
Semi-VOC	EPA Method	Y				INF							
Method 8270C	8270C			Y		EFF							
Maral.	See note 3	D/Y				INF	INF	INF(T)	INF(T)	INF(T)			
Metals	below			D/Y		EFF	EFF	EFF(T)	EFF(T)	EFF(T)			
Low level mercury (if	EPA Method	D/Y					INF						
needed)	1631A			D/Y			EFF				1		
Cyanide Total	Total CN - EPA Method	D/Y				INF	INF						
Cyaniac Total	335.2			D/Y		EFF	EFF						

						4th Q	uarter of Ope	eration	5th Q	uarter of Ope	eration
Type of Sample	Analytical Method	INF-001	MID-001	EFF-001	RSW-001U/ RSW-001D	Oct-11	Nov-11	Dec-11	Jan-12	Feb-12	Mar-12
		Q						INF			INF
PAH method 8310	EPA Method 8310			Q				EFF			EFF
					V						
		Y									
EDB method 504	EPA Method 504			Y							
					V	-					
		Y				INF	INF	INF	INF	INF	INF
VOC method	EPA Method		Y			MID	MID	MID	MID	MID	MID
8260b	8260B			Y		EFF	EFF	EFF	EFF	EFF	EFF
					V						
		Y									
Fuel Oxygenates TAME, DIPE, ETBE, TBA, Ethanol, Methanol (with	EPA Method 8260B			Q (annual if not reported above 5 ug/l in first sample)	-1-	-1		EFF			EFF
VOC analyses)					V						
Semi-VOC	EPA Method	Y									
Method 8270C	8270C			Y							
Metals	See note 3	D/Y									
Wietais	below			D/Y							
Low level mercury (if	EPA Method	D/Y									
needed)	1631A			D/Y							
Cyanide Total	Total CN - EPA Method	D/Y									
- y	335.2			D/Y							

						1st Qua	arter of ation	2nd Q	uarter of Ope	eration	3rd Q	uarter of Ope	eration
Type of Sample	Analytical Method	INF-001	MID-001	EFF-001	RSW-001U/ RSW-001D	Star	tup	Apr-11	May-11	Jun-11	Jul-11	Aug-11	Sep-11
						Day 1	Day 5						
Flow Rate (gpm and gpd)	Field Measurement			con-tinuous									
Turbidity	EPA Method 180.1			D/Q/Y		EFF	EFF	EFF	EFF	EFF	EFF	EFF	EFF
		D/M/Q/Y				INF	INF	INF	INF	INF	INF	INF	INF
рН	EPA Method 150.1			D/M/Q/Y		EFF	EFF	EFF	EFF	EFF	EFF	EFF	EFF
					V								
Dissolved Oxygen (mg/l)	EPA Method 360.1				V								
Total Dissolved Solids				D/M		EFF	EFF	EFF	EFF	EFF	EFF	EFF	EFF
	Field	D				INF	INF						
Temperature	Measurement			D/M/Q/Y		EFF	EFF	EFF	EFF	EFF	EFF	EFF	EFF
		D				INF	INF						
Electrical Conductivity	EPA Method 120.1			D/M/Q/Y		EFF	EFF	EFF	EFF	EFF	EFF	EFF	EFF
Hardness as CaCO ₃	SM				T			EFF(T)	EFF(T)	EFF(T)			
Salinity	SM				T			EFF(T)	EFF(T)	EFF(T)			

						4th Q	uarter of Ope	eration	5th Q	uarter of Ope	eration
Type of Sample	Analytical Method	INF-001	MID-001	EFF-001	RSW-001U/ RSW-001D	Oct-11	Nov-11	Dec-11	Jan-12	Feb-12	Mar-12
Flow Rate (gpm and gpd)	Field Measurement			con-tinuous							
Turbidity	EPA Method 180.1			D/Q/Y		EFF	EFF	EFF	EFF	EFF	EFF
		D/M/Q/Y				INF	INF	INF	INF	INF	INF
рН	EPA Method 150.1			D/M/Q/Y		EFF	EFF	EFF	EFF	EFF	EFF
					V						
Dissolved Oxygen (mg/l)	EPA Method 360.1				V	1					
Total Dissolved Solids				D/M		EFF	EFF	EFF	EFF	EFF	EFF
	Field	D									
Temperature	Measurement			D/M/Q/Y		EFF	EFF	EFF	EFF	EFF	EFF
		D									
Electrical Conductivity	EPA Method 120.1			D/M/Q/Y		EFF	EFF	EFF	EFF	EFF	EFF
Hardness as CaCO ₃	SM				T						
Salinity	SM				T						

Table 2-1

Metals	Total (unfiltered) pp13 metals - Mercury, cadmium, silver, nickel, thallium, zinc, arsenic selenium, antimony, berylium, total chorium, copper, and lead.
Inorganic	Must report with at most these limits (in micrograms per liter [ug/L]): Mercury - 0.002; Cadmium and Silver - 0.25; Nickel, Tahllium and Zinc - 1.0; Arsenic and Selenium - 2.0; Cyanide - 1.0; Antimony,
reporting limits:	Beryllium, Ttoal Chromium, Copper and Lead (0.5).
D	Once during the first and fifth day of startup
M	Once each month
Q	Once during the first week of startup, quarterly thereafter, unless below 0.5 u/l, then annual thereafter
Y	Once during the first week of startup, annually thereafter
D/W/M	Once during the first and fifth day of startup, weekly for the remainder of the month, monthly thereafter
D/M	Once during the first and fifth day of startup, monthly thereafter
D/Q	Once during the first and fifth day of startup, quarterly thereafter
D/Y	Once during the first and fifth day of startup, annually thereafter
Q/Y	Quarterly for the first year of operation, annually thereafter
D/Q/Y	Once during the first and fifth day of startup, monthly for the first year of operation, annually thereafter
D/M/Q/Y	Once during the first and fifth day of startup, monthly for the first year of operation, quarterly for the second year, and annually thereafter. In case of pH analyses, this monitoring requirement is only for facilities with treatment processes that would cause no pH variances in the effluent. If any chemical used in the treatment process may cause pH variances in the effluent, the frequency of pH monitoring in the effluent shall be increased to twice per week for the first month of operation and weekly thereafter if pH monitoring data for the first month of operation demonstrate compliance with pH effluent limits.
V T EFF-T/INF-T	Sampling should be performed within 24 hours after an effluent limit violation is confirmed in EFF-001 Sampling should be performed when Cadmium, chromium(total), Copper, Lead, Nickel, Silver or Zinc triggers are exceeded. Sampling added for trigger evaluation during Q22011. Analyze metals for total cadmium and total copper only.

TABLE 3-1 FLOW AND OPERATIONS SUMMARY

Startup and First Quarter 2011

Greenway Phase II NDPES Dewatering Treatment System Hollis Street, Emeryville, California

		Elapsed Time	System Average Flow Rate ¹	System Cumulative Volume ²	COMMENTS
Date	Time	(days)	(gpm)	(gallons)	-
3/3/2011	725	0.00	0.0		System begins discharging to storm drain
3/3/2011	910	0.08	66.0	7,910	Batch Tank pumped down at 1048
3/3/2011	1535	0.25	14.1	12,990	
3/4/2011	835	0.71	5.2	18,310	
3/7/2011	730	2.96	5.4	41,520	
3/8/2011	730	1.00	2.3	44,850	
3/9/2011	930	1.08	3.1	49,680	
3/11/2011	1330	2.17	4.8	64,590	Repaired leak in 3 inch hose at batch tank, Backwashed sand filters.
3/14/2011	800	2.79	5.2	85,670	Changed out 3 inch hose
3/24/2011	800	10.00	7.9	200,120	Manual backwash sand and carbon vessels
3/28/2011	800	4.00	18.5	306,500	Manual backwash sand and carbon vessels

Summary

			Julilliai y		
	Elapsed Time	Average Flow	Cumulative	Daily Average	Total Volume Pumped During Period
	in Period	Rate ¹	Volume ²	Flow Rate	Total Volume Pumped During Period
	(days)	(gpm)	(gallons)	(gpd)	(gallons)
Startup Period	8.3	5.4	64,590	7,829	64,590
First Quarter 2011 ⁽³⁾	25.0	8.5	306,500	12,240	306,500
	Permitted Flow	100			

Notes:

 $^{^{1}}$ System average flow rate is calculated as the change in system cumulative volume (from the previous measurement to the current measurement) divided by the elapsed time in minutes.

² Cumulative volume since startup

³ Includes Startup Period data

SUMMARY OF

GROUNDWATER TREATMENT SYSTEM PARAMETERS

Startup and First Quarter 2011

Greenway Phase II NPDES Dewatering Treatment System Hollis Street, Emeryville, California

Sample	Date	Туре	рН		Electrical Cond	uctivity	Turbidity	,	TDS	Ammonia as N (1)	Alkalinity as CaCO ₃ ⁽¹⁾	Hardness as CaCO ₃ ⁽¹⁾	Dissolved Oxygen ⁽¹⁾	96-hour Static Fish Bioassay ⁽²⁾
Location		(field/lab)	(S.U.)	%	(µmhos/cm)	%	(NTUs)	%	(mg/l)	(mg/l)	(mg/l)	(mg/l)	(mg/l)	survival)
		EPA Method	field/ 9040C	differ- ence	field/SM2510B	differ- ence	field/SM2130B	differ- ence	SM2540C		ЕРА М	ethod EPA/82	1/R-02/012	
La	boratory Re	eporting Limit	1		10		0.1		13					
	2/28/2011	field	7.52		1,270		69.7							
INIE 001	2/20/2011	lab			1,100									
INF-001	3/7/2011	field	7.4		1,460		39.6							
	3/1/2011	lab	7.37										1	
		field	7.82/7.91/8.05	1.6	1230/1250/1240	0.8	28.7/29.1/29.3	1.1						
	2/28/2011	lab	8.40		1,100				730					
EFF-001	3/7/2011	field	7.29/7.27/7.34	0.5	1430/1450/1440	0.7	18.2/17.6/17.4	2.4						
	3/1/2011	lab	7.81		1,300		2.7		800					
	Effluent Limitati		6.5-8.5	See Note 6		See Note 6		See Note 6						> 70 percent ³ > 90 percent ⁴
	Trigger Levels						5							

Notes:

NTUs = nephelometric turbidity ur

°C= degrees Celsius

S.U. = standard units

ımhos/cm= micromhos per centimeter

(1) Analyzed as part of Fish Bioassay

(2) Test species - rainbow trout

(3) Greater than 70 percent (single test)

(4) Greater than 90 percent (3-sample moving average)

(6) Sample when deviation from average is less than 5 percent

TABLE 3-3 SUMMARY OF CHEMICAL ANALYSES FOR ORGANIC POLLUTANTS

Startup and First Quarter 2011

Compound	l Number in Permit	1	2	3	4	5	6	7	8	9	10
				·		VO	Cs				
			Carbon	1,1-	1,2-	1,1-			Tetra-		cis-1,2-
Sample			Tetra-	Dichloro-	Dichoro-	Dichloro-	Ethyl	Methyl-ene	chloro-		Dichloro-
Location/	and the second	Benzene	chloride	ethane	ethane	ethene	benzene	Chloride	ethene	Toluene	ethene
ID	Date Sampled	(µg/l)	(µg/l)	(µg/l)	(µg/l)	(µg/l)	$(\mu g/l)$	(µg/l)	(µg/l)	(µg/l)	(µg/l)
	EPA Method	8260B	8260B	8260B	8260B	8260B	8260B	8260B	8260B	8260B	8260B
Labora	tory Reporting Limit	0.5	0.5	0.5	0.5	0.5	0.5	5.0	0.5	0.5	0.5
Ме	thod Detection Limit	0.075	0.072	0.067	0.077	0.058	0.070	1.0	0.065	0.17	0.071
INF-001	2/28/2011	ND	ND	ND	ND	ND	ND	ND	ND	0.25 J	ND
INF-001	3/7/2011	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
MID 001	2/28/2011										
MID-001	3/7/2011	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
EEE 001	2/28/2011	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
EFF-001	3/7/2011	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Effluent Lii	mitations	1.0	0.5	5.0	5.0	0.11	5.0	5.0	1.6	5.0	5.0

TABLE 3-3 SUMMARY OF CHEMICAL ANALYSES FOR ORGANIC POLLUTANTS

Startup and First Quarter 2011

Compound	l Number in Permit	11	12	13	14	15	16	17	18	19	20
					VOCs				T1	PH	
			1,1,1-	1,1,2-							
Sample		trans-1,2-	Trichloro-	Trichloro-	Trichloro-	Vinyl	Total		TPH as	TEPH as	
Location/		Dichloro-ethene	ethane	ethane	ethene	Chloride	Xylenes	MTBE	Gasoline	Diesel Fuel	EDB
ID	Date Sampled	$(\mu g/l)$	(µg/l)	$(\mu g/l)$	(µg/l)	(µg/l)	(µg/l)	$(\mu g/l)$	(µg/l)	(µg/l)	$(\mu g/l)$
	EPA Method	8260B	8260B	8260B	8260B	8260B	8260B	8260B	8260B	8015M	504.1
Labora	tory Reporting Limit	0.5	0.5	0.5	0.5	0.5	1	0.5	51	52	0.02
Ме	thod Detection Limit	0.070	0.055	0.11	0.059	0.05	0.49	0.069	24	25	0.0029
INF-001	2/28/2011	ND	ND	ND	0.083 J	0.1 J	ND	ND	ND	ND	
INF-001	3/7/2011	ND	ND	ND	0.11 J	ND	ND	ND	ND	24 J B	ND
MID 001	2/28/2011										
MID-001	3/7/2011	ND	ND	ND	ND	ND	ND	ND	ND		
EEE 001	2/28/2011	ND	ND	ND	ND	ND	ND	ND	ND	ND(<24)	
EFF-001	3/7/2011	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Effluent Lii	mitations	5.0	5.0	1.2	5.0	0.5	5.0	5.0 ⁽²⁾	50	50	< 0.05

SUMMARY OF CHEMICAL ANALYSES FOR ORGANIC POLLUTANTS

Startup and First Quarter 2011

Greenway Phase II NPDES Dewatering Treatment System Hollis Street, Emeryville, California

Notes:

 $\mu g/l = micrograms per liter$

VOCs = volatile organic compounds, analyzed by EPA Method 8260B

SVOCs = semi-volatile organic compounds, analyzed by EPA method 8270C

EDB = Ethylene Dibromide

gasoline

TPH as = total petroleum hydrocarbons C6-C12

TEPH as (EPA Method 3630C) = total extractable petroleum hydrocarbons, diesel - C10-C24, motor oil - C24-C36, precleaned with Silica Gel cleanup (EPA Method 3630C)

Diesel Fuel

MDL = Method Detection Limit (range shown if MDL varies between analyses)

-- = Not analyzed

ND = not detected above method detection limit

J = J-flag by laboratory, indicating estimated concentration below reporting limit but above method detection limit

Lab does not report 1,1,2 trichlorotrifluoroethane

Based on Toxics pollutants defined in the Fuel General NPDES Permit No. CAG912002, Section IV.A.1,

Table 2: Effluent Limitations for Toxics Pollutants.

SUMMARY OF CHEMICAL ANALYSES FOR ORGANIC TRIGGER COMPOUNDS $^{\left(1\right)\left(2\right)}$

Startup and First Quarter 2011

Greenway Phase II NPDES Dewatering Treatment System

Hollis Street, Emeryville, California

			VOCs			Volatile Fu	el Additiv	es		sVC	OCs	PAI	Hs
			Carbon	1,2,4- Trimethyl-					TEPH as Motor Oil	Bis (2- ethylhexyl)	Other		Other
		Acetone	Disulfide	benzene	TBA	DIPE	ETBE	TAME		phthalate	sVOCs	Fluorene	PAHs
Sample Location	Date Sampled	$(\mu g/l)$	$(\mu g/l)$	(µg/l)	$(\mu g/l)$	$(\mu g/l)$	(µg/l)	(µg/l)	(µg/l)	(µg/l)	$(\mu g/l)$	(µg/l)	(µg/l)
	EPA Method	8260B	8260B	8260B	8260B	8260B	8260B	8260B	8015M	8270C	8270C	8310	8310
Laboratory	Reporting Limit	50	5	0.5	4	0.5	0.5	0.5	100	10	Varies	0.1	0.1
Method	l Detection Limit	3.7	0.078	0.045	1.9	0.05	0.098	0.071	37	1.5	Varies	0.032	Varies
INF-001	2/28/2011	11 J	ND	ND	ND	1.5	ND	ND	51 J B	3.7 J	ND		
1141 -001	3/7/2011	ND	ND	ND	ND	0.88	ND	ND	ND			0.050 J	ND
MID-001	2/28/2011					1				ŀ			
WIID-001	3/7/2011	ND	ND	0.070 J B	ND	ND	ND	ND					
EFF-001	2/28/2011	ND	ND	0.094 J B	ND	ND	ND	ND	41 J B	ND	ND		
Li i -001	3/7/2011	ND	0.37 J	ND	ND	ND	ND	ND	50 J B			ND	ND
	Trigger Levels	5.0	5.0	5.0	5.0	5.0	5.0	5.0	50	5	5	5	Varies
Trigger Evalu	ation Warranted					X			X				

SUMMARY OF CHEMICAL ANALYSES FOR ORGANIC TRIGGER COMPOUNDS (1)(2)

Startup and First Quarter 2011

Greenway Phase II NPDES Dewatering System

Hollis Street, Emeryville, California

Notes:

- $^{\left(1\right)}$ Only trigger compounds with concentrations greater than MDL are listed in the table.
- (2) Based on Organic Trigger Compounds defined in the Fuel General NPDES Permit No. CAG912002, Section VI.C.6, Table 3: Trigger Compounds or Constituents.
- MDL = method detection limit
- TEPH = total extractable petroleum hydrocarbons, diesel C10-C24, motor oil C24-C36, precleaned with Silica Gel cleanup (EPA Method 3630C)
- $\mu g/l = micrograms per liter$
- VOC = volatile organic compound
- SVOC = semi-volatile organic compound
- ND = not detected above method detection limit
- J = J-flag by laboratory, indicating concentration is below reporting limit but above method detection limit
- B = Compound was found in the blank and sample.
- ↑ Trigger level exceeded in at least one sample
- b Lab narrative states high response in CCD
- TBA = tert-butyl alcohol
- DIPE = Isopropyl Ether
- ETBE = Ethyl tert-Butyl Ether
- TAME = Methyl tert-Amyl Ether
- PAHs = polynuclear aromatic hydrocarbons

SUMMARY OF CHEMICAL ANALYSES FOR METALS, METALLOIDS AND CYANIDE

Startup and First Quarter 2011

Greenway Phase II NPDES Dewatering Treatment System

Hollis Street, Emeryville, California

Sample Location/ ID	Date Sampled	Antimony (μg/l)	Arsenic (µg/l)	Berylliu m (µg/l)	Cadmiu m (µg/l)	Chromium (µg/l)	Copper (µg/l)	Lead (µg/l)	Mercury (µg/l)	Low-Level Mercury (µg/l)	Nickel (µg/l)	Selenium (µg/l)	Silver (µg/l)	Thallium (µg/l)	Zinc (µg/l)	Cyanide (µg/l)
EPA A	nalytical Method	6020	6020	6020	6020	6020	6020	6020	7470A	1631	6020	6020	6020	6020	6020	SM4500 CN- E
Laborator	y Reporting Limit	2.0	1.0	0.5	1.0	2.0	2.0	1.0	0.2	0.0005	2.0	2.0	1.0	1.0	20	3
Metho	d Detection Limit	0.3	0.9	0.1	0.1	0.9	0.5	0.2	0.1	0.00012	0.5	0.5	0.1	0.2	4	2.2
INF-001	2/28/2011	0.78 J	1.5	ND	0.11 J	7.0	33	3.0	ND		11	1.6 J	ND	ND	50	ND (<17)
1111-001	3/7/2011	0.56 J	ND	ND	0.21 J	1.7 J	5.4	0.41 J	ND	0.0109 J	7.1	ND	ND	ND	10 J	ND
EFF-001	2/28/2011	0.59 J	ND	ND	ND	2.1	6.2	1.1	ND		3.4	1.3 J	ND	ND	4.6 J	ND (<17)
EFT-001	3/7/2011	0.58 J	ND	ND	ND	ND	1.3 J	ND	ND	0.0066 J	3.0	ND	ND	ND	ND	ND
Trigger C	Concentration	6	10	1	0.07	11	3.1	2.0	0.025	0.025	8.2	5.0	1.9	0.1	35	1.0
Trigge	er Evaluation Warranted				x		x									

Notes:

 $\mu g/l = micrograms per liter$

ND = not detected above method detection limit

J = J-flag by laboratory, indicating concentration is below reporting limit but above method detection limit

= Trigger

X evaluation

warranted

TABLE 3-6 SUMMARY OF MASS REMOVAL - ORGANIC COMPOUNDS Startup and First Quarter 2011 Greenway Phase II NPDES Treatment System Hollis Street, Emeryville, California

Period	DIPE	A Toluen		ge Influer TCE	nt Co	oncentratio Vinyl Chlorid		Total VOCs	Volume of Water Treated	Mass of VOCs Removed	Cumulative VOC Mass Removed
	(µg/l)	(µg/l)		(µg/l)		(µg/l)		$(\mu g/l)$	(gallons)	(pounds)	(pounds)
System Startup	1.2	0.17	J	0.10	J	0.06	J	1.5	64,590	0.0008	0.0008
First Quarter 2011	1.2	0.2	J	0.1	J	0.1	J	1.5	306,500	0.0039	0.0039

Notes:

For calculation of averages, when compound is detected above method detection limits in at least one sample in the quarter, then, for samples during the quarter in which the concentration is below the method detection limit, one half of MDL is used. If the compound is not detected in any samples during the quarter, then compound is not included in the calculation of mass removed.



FIGURES



Site Location Map Greenway Phase II NPDES Treatment System 5812 Hollis Street, Emeryville, CA

Figure 1-1

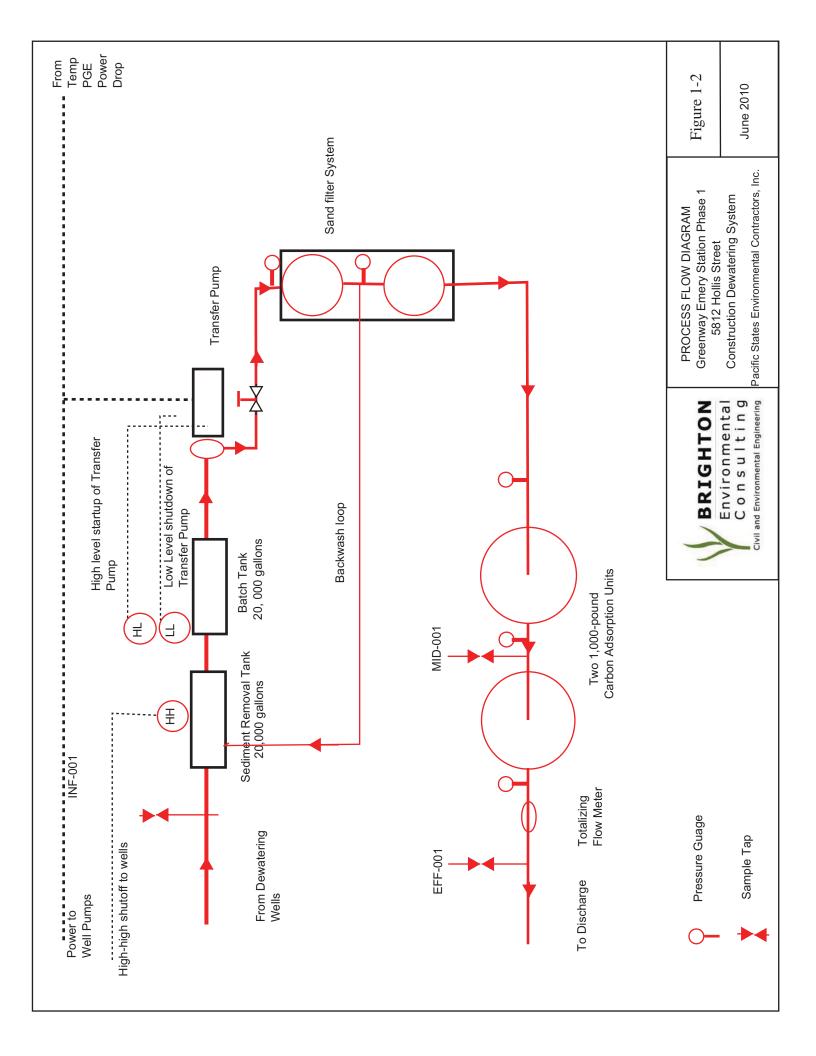
June 2010

Approximate North

1200

Approximate scale (feet)







APPENDIX A QA/QC DOCUMENTATION

TABLE A-1 SUMMARY OF SAMPLE COLLECTION METHODS

Analyses	Method	COMMENTS
Turbidity	SM 2130B	Each sample collected in unpreserved 500 ml plastic bottle
Total Dissolved Solids	SM 2450C	Each sample collected in unpreserved 500 ml plastic bottle
Bioassay 96 hour static renewal survival	EPA-821-R-02-012	Each sample collected in unpreserved 5 gallon plastic container
рН	EPA Method 9040C	Each sample collected in unpreserved 500 ml plastic bottle
Electrical Conductivity	SM 2510B	Each sample collected in unpreserved 500 ml plastic bottle
Soluble Metals	EPA Method 6000 series	Each sample collected in two 500 ml plastic bottles unpreserved (filtered and preserved with nitric acid at lab within 24 hours)
Total Metals	EPA Method 6000 series	Each sample collected unfiltered in two 500 ml plastic bottles with nitric acid
Low level Mercury	EPA Method 1631A	One 500 ml with HCl, special clean gloves/dirty gloves handling techniques in field.
Cyanide Total	Kelada-01	500 ml with NaOH
Ethylene Dibromide (EDB)	EPA Method 504	Each sample collected in three 40-milliliter (ml) glass vials closed with screw caps with TeflonTM-lined septa, containing hydrochloric acid (HCl) for sample preservation.
Fuel Oxygenates TAME, DIPE, ETBE,		Each sample collected in three 40-milliliter (ml) glass vials
TBA, Ethanol, Methanol Volatile Organic Compounds	EPA Method 8260B	closed with screw caps with Teflon TM -lined septa, containing hydrochloric acid (HCl) for sample preservation.
VOC method 8021 (includes BTEX and MTBE)	BTEX -EPA Method 8021	Each sample collected in three 40-milliliter (ml) glass vials closed with screw caps with TeflonTM-lined septa, containing hydrochloric acid (HCl) for sample preservation.
Semi-volatile organic componds	EPA Method 8270C	Each sample collected in two HCL preserved 1 liter amber glass bottle
PAH method 8310	EPA Method 8310	Each sample collected in two HCL preserved 1 liter amber glass bottle
Total petroleum hydrocarbons as gasoline	EPA Method 8015 Mod	Each sample collected in three 40-milliliter (ml) glass vials closed with screw caps with Teflon TM -lined septa, containing hydrochloric acid (HCl) for sample preservation.
TPHd (8015M) with silica gel cleanup (diesel and motor oil)	EPA Method 8015 Mod with silica gel cleanup	Each sample collected in one HCL preserved 1 liter amber glass bottle

${\it TABLE~A-2} \\ {\it SUMMARY~OF~ANALYTICAL~METHODS~AND~LABORATORY~QA/QC} \\$

Analyte	Analysis Method	Analysis Sample Date	Performed by Discharger	Performed by Analytical Laboratory	Laboratory Name	Is lab State-certified in these methods?	Analyses pertormed according to standard methods?	Sample holding times met?	Analytical results reported when above the method detection or reporting limit?	Reporting limits adequate to determine compliance?	QA/QC analyses run consistent with analytical methods?	QA/QC results meet all acceptance criteria?
Flow rate	field	Weekly	X									
Turbidity	SM 2130B/field	2/28/2011	X	X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
Turbianty	SWI 2130B/Held	3/7/2011	X	X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
Total Dissolved	SM 2540C	2/28/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
Solids	SWI 2540C	3/7/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
	field	Monthly	X			yes	yes	yes	yes		yes	yes
pН	EPA 9040C	2/28/2011	X	X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
	LI A 7040C	3/7/2011	X	X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
Standard Observations	field	Monthly	X									
Temperature	field	Monthly	X									
Antimony	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Anumony	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Arsenic	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Arsenie	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Beryllium	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Dorymani	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Cadmium	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Copper	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Соррег	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes

TABLE A-2 SUMMARY OF ANALYTICAL METHODS AND LABORATORY QA/QC

Analyte	Analysis Method	Analysis Sample Date	Performed by Discharger	Performed by Analytical Laboratory	Laboratory Name	Is lab State-certified in these methods?	Analyses performed according to standard methods?	Sample holding times met?	Analytical results reported when above the method detection or reporting limit?	Reporting limits adequate to determine compliance?	QA/QC analyses run consistent with analytical methods?	QA/QC results meet all acceptance criteria?
Lead	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Doug	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Mercury	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
	7470	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Low Level Mercury	EPA Method 1631	3/7/2011		X	TANC ⁽³⁾	yes	yes	yes	yes	yes	yes	yes
Nickel	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
TVICKCI	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Selenium	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Scientini	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Silver	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Sirver	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Thallium	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Zinc	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Zinc	6010B	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Cyanide	EPA Method	2/28/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
Cyamac	335.2	3/7/2011		X	TAI ⁽¹⁾	yes	yes	yes	yes	yes	yes	yes
VOCs	EPA Method	2/28/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	no	yes	yes
8021B or 8260B	3/7/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes	

TABLE A-2 SUMMARY OF ANALYTICAL METHODS AND LABORATORY QA/QC

Analyte	Analysis Method	Analysis Sample Date	Performed by Discharger	Performed by Analytical Laboratory	Laboratory Name	Is lab State-certified in these methods?	Analyses performed according to standard methods?	Sample holding times met?	Analytical results reported when above the method detection or reporting limit?	Reporting limits adequate to determine compliance?	QA/QC analyses run consistent with analytical	QA/QC results meet all acceptance criteria?
TPH as gasoline	EPA 8015M	2/28/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
1111 us gusonne		3/7/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
TPH as diesel	EPA 8015M	2/28/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
11 11 ds diesei	L171 0015W1	3/7/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
TEPH as motor	EPA 8015M	2/28/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	no(5)	yes	yes
oil	Li ii oo i siyi	3/7/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	no(5)	yes	yes
EDB	EPA Method 504	3/7/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	yes
SVOCs	EPA Method 8270C	2/28/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	yes	yes	Method 8270C: The laboratory control sample and the laboratory control sample duplicate (LCS/LCSD) for batch #86975 exceeded control limits for the following analyte(s): Benzoic acid. Benzoic acid has been identified as a poor performing analyte when analyzed using this method; therefore, re-extraction/re-analysis wa not performed. Batch precision also exceeded control limits for these analyte(s). These results have been reported and qualified. Method 8270C: A full list spike was utilized for this method. Due to the large number of spiked analytes, there is a high probability that one or more analytes will recover outside acceptance limits. The laboratory's SOP allows for three (03) analytes to recover outside criteria for this method hen a full list spike is utilized. The LCSD associated with batch #86975 had one (Dimethylphthalate) analyte outside control limits; therefore, re-extraction/re-analysis as not performed. These results have been reported and qualified.
PAHs	EPA Method 8310	3/7/2011		X	TASF ⁽²⁾	yes	yes	yes	yes	no ⁽⁵⁾	yes	yes

TABLE A-2 SUMMARY OF ANALYTICAL METHODS AND LABORATORY QA/QC

Greenway Phase II NPDES Treatment System Hollis Street, Emeryville, California

	Analysis	nalysis Sample Date	Performed by Anal	ormed by ytical Laboratory	lab State-certified in these sthods? natyses performed cording to standard	mple holding times met?	nalytical results reported nen above the method tection or reporting limit?	Forting limits adequate to termine compliance? A/CC analyses run nsistent with analytical schools?	A/QC results meet all ceptance criteria?	
Analyte	Method	Ana	1 - 1	ratory Name	Is Is Is met Ang acc	San	Ana whe dete	Reg detre	QA	

Notes:

-- = not applicable

(3) TestAmerica North Canton, 4101 Shuffel Street, NW, North Canton, OH 44720

NA- not analyzed

(4) Laboratory Reporting limit exceeds trigger concentration, MDL is below trigger concentrations

(1) TestAmerica Irvine, 17461 Derian Avenue. Suite 100, Irvine, CA 92614

⁽⁵⁾Laboratory reporting limit and MDL exceed trigger level.

⁽²⁾ TestAmerica San Francisco, 1220 Quarry Lane, Pleasanton, CA 94566



APPENDIX B STANDARD OBSERVATIONS

TABLE B-1 STANDARD OBSERVATIONS OF STORM DRAIN INLET

Greenway Phase II NPDES Treatment System Emeryville, California

Standard Observations for Treatement System Date	Odor	Weather Condition	Deposits, discoloration and or plugging in system that could adversely affect system reliability		COMMENTS
2/28/2011	no	clear	no	na	
3/3/2011	no	cloudy	no	na	
3/7/2011	no	cloudy	no	na	