P & D Environmental

A Division of Paul H. King, Inc. 4020 Panama Court Oakland, CA 94611 (510) 658-6916

April 19, 2004 Report 0055.R18

Mr. Manmohan Chopra 4216 Warbler Loop Fremont, CA 94555

SUBJECT: GROUNDWATER MONITORING AND SAMPLING REPORT

Former Haber Oil Station 1401 Grand Avenue San Leandro, California

Dear Mr. Chopra:

P&D Environmental (P&D) is pleased to present this report documenting the results of the most recent monitoring and sampling of the eight-groundwater monitoring wells at or near the subject site. This work was performed in accordance with P&D's proposal 022698.P1 dated February 26, 1998. All of the wells were monitored and sampled on November 25, 2003. A Site Location Map (Figure 1) and Site Plan (Figure 2) are attached with this report.

BACKGROUND

The site is presently used as an active gasoline station. It is P&D's understanding that on April 24, 1991 Aegis Environmental, Inc. (Aegis) personnel drilled four soil borings, designated as B-1 through B-4, to a vertical depth of approximately 40 feet at the site. The locations of the borings are shown on Figure 2. A total of nine soil samples collected from the boreholes were analyzed for total petroleum hydrocarbons as gasoline (TPH-G); benzene, toluene, ethylbenzene, and total xylenes (BTEX) by EPA Method 8260; and for total lead by EPA Method 7420. TPH-G concentrations ranged from below detection limit to 66 milligram per kilogram (mg/kg). Benzene concentrations ranged from not detected to 0.94 mg/kg. Total lead concentrations ranged from not detected to 3 mg/kg. Documentation of the subsurface investigation and results are presented in a report prepared by Aegis titled, "Soil Boring Results Report," dated June 10, 1991.

It is P&D's understanding that on April 14, 1992 Aegis personnel returned to the site to drill three slant borings, designated as B5 through B7, to a total vertical depth of approximately 49 feet at the site. The borings were drilled at an angle of approximately 26 to 28 degrees to collect samples from beneath the underground storage tanks. The locations of the borings are shown on Figure 2. A total of twenty-two soil samples were analyzed for TPH-G using EPA Method 5030, and for BTEX using EPA Method 8240. In addition, one of the samples was analyzed for total lead using EPA Method 7420, and several of the soil samples were analyzed for soluble lead using the California Waste Extraction Test. TPH-G concentrations ranged from not detected to 4,000 mg/kg. Benzene, concentrations ranged from not detected to 11 mg/kg. Total lead was not detected, and soluble lead concentrations ranged from not detected to 0.061 mg/kg. Documentation of the subsurface investigation and results are presented in a report prepared by Aegis titled, "Initial Subsurface Investigation Results Report," dated June 22, 1992.

It is P&D's understanding that between September 15 and 18, 1992 Aegis personnel returned to the site to install five groundwater monitoring wells, designated as MW1 through MW5. The

wells were drilled to total depths of between 50 and 55 feet, and were constructed using four-inch diameter PVC pipe. Wells MW1 and MW2 were constructed with perforated casing between the depths of approximately 15 and 55 feet. Wells MW3, MW4 and MW5 were constructed with perforated casing between the depths of approximately 35 and 55 feet. Groundwater was reported as first encountered at a depth of 42 feet. The well locations are shown in Figure 2.

A total of thirty-one soil samples were analyzed for TPH-G using EPA Method 5030/8015; and for BTEX using EPA Method 8020. In addition, three soil samples containing TPH-G were analyzed for total metals concentrations of cadmium, chromium, lead, and zinc using EPA Method 6010 and 7421. One soil sample was collected from each borehole from below the airwater interface and analyzed for petrophysical properties, including saturated permeability and grain size distribution.

TPH-G concentrations ranged from not detected to 39 mg/kg. Benzene concentrations ranged from not detected to 0.27 mg/kg. The total metals concentrations were all less than 10 times their respective STLC values. The subsurface materials encountered in the borings indicate that soil types vary across the site, but generally consist of silty clay, silt, clayey silt and sandy silt from the surface to a depth of between 30 and 35 feet. Below the depth of 30 to 35 feet, layers of sand and sandy silt were reported to have been encountered.

It is P&D's understanding that on September 29, 1992 Aegis personnel collected groundwater samples from wells MW1, MW2, MW4 and MW5 at the site. A sample was not collected from well MW3 due to the reported presence of 0.02 feet of floating hydrocarbons. The measured depth to water ranged from approximately 41.5 to 44.5 feet. The samples were analyzed for TPH-G using EPA Method 5030/8015; and for BTEX using EPA Method 8020. TPH-G concentrations ranged from 0.06 to 20 mg/kg, and benzene concentrations ranged from 0.16 to 10 mg/kg. Based upon the water level measurements in the wells, the groundwater flow direction was reported to be to the northwest. The water level measurements are summarized in Table 1. The analytical results are summarized in Table 2.

It is P&D's understanding that on October 7, 1992 Aegis personnel performed rising head slug tests wells MW1, MW2, and MW4 to estimate the saturated hydraulic conductivity at the site. In addition, two short-term soil vapor extraction tests were performed on wells MW1 and MW2. Wells MW3, MW4, and MW5 were used as vacuum influence monitoring points. Documentation of the monitoring well groundwater sample collection, slug test and vapor extraction tests are presented in a report prepared by Aegis titled, "Problem Assessment Report," dated December 16, 1992.

On February 18, 1994 P&D personnel monitored the five-groundwater monitoring wells at the site for depth to water and the presence of free product or sheen. The depth to water was measured using an electric water level indicator, and the presence of free product and sheen was evaluated using a transparent bailer. The measured depth to water in the wells ranged from approximately 39.8 to 42.9 feet. No evidence of free product or sheen was detected in any of the wells. Based on the measured depth to water in the wells, the groundwater flow direction was calculated to be to the north with a gradient of 0.054. In a letter dated October 19, 1995 Mr. Scott Seery of the Alameda County Department of Environmental Health requested that all of the

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onsite and offsite wells be monitored and sampled for the quarterly monitoring and sampling program. The measured depth to water in the wells is presented in Table 1.

On June 15 and 16, 1995 P&D installed three offsite monitoring wells, designated as MW6 through MW8. The locations of the wells are shown on Figure 2. Documentation of the well installation and sample results is presented in P&D's report 0055.R5 dated August 23, 1995.

The underground storage tanks at the subject site were replaced in the first half of 1997. Following removal of the tanks, excavation of soil was performed in the area surrounding well MW1. As a result of the excavation activities, the elevation at the top of well MW1 was altered. The present elevation for the top of well MW1 is unknown.

In January 2003 Ms. Eva Chu of the Alameda County Department of Environmental Health (ACDEH) requested that the wells be analyzed for fuel oxygenates using EPA Method 8260. In a letter dated June 20, 2003 Ms. Chu requested that the analysis be continued.

FIELD ACTIVITIES

On November 25, 2003 all eight of the wells in the groundwater monitoring network for the site were monitored, purged, and sampled by P&D personnel. The wells were monitored for depth to water and the presence of free product or sheen. Depth to water was measured to the nearest 0.01 foot using an electric water level indicator. The presence of sheen was evaluated using a transparent bailer. No free product or sheen was observed in any of the wells. Depth to water level measurements and monitoring well groundwater surface elevations are presented in Table 1.

Prior to sampling, the wells were purged of a minimum of three casing volumes of water. During purging operations, the field parameters of electrical conductivity, temperature and pH were monitored. Once the field parameters were observed to stabilize, and a minimum of three casing volumes had been purged, water samples were collected using a clean Teflon bailer. The water samples were transferred to 40-milliliter glass Volatile Organic Analysis (VOA) vials, which were sealed with Teflon-lined screw caps. The VOA vials were overturned and tapped to assure that no air bubbles were present.

The VOA vials were then transferred to a cooler with ice, and later were transported to McCampbell Analytical, Inc. in Pacheco, California. McCampbell Analytical, Inc. is a State-certified hazardous waste testing laboratory. Chain of custody documentation accompanied the samples to the laboratory. Records of the field parameters measured during well purging are attached with this report.

HYDROGEOLOGY

The subsurface materials encountered in the borings drilled by Aegis indicate that soil types vary across the site, but generally consist of silty clay, silt, clayey silt and sandy silt from the surface to a depth of between 30 and 35 feet. Below the depth of 30 to 35 feet, layers of sand and sandy silt were reported to have been encountered. Groundwater has historically been encountered at the site at depths ranging from approximately 40 to 45 feet below grade.

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Based upon the regional groundwater flow direction identified by Woodward-Clyde Consultants in a report titled, "Hydrogeology of Central San Leandro and Remedial Investigation of Regional Groundwater Contamination - San Leandro Plume - San Leandro, California - Volume I," prepared for the California Environmental Protection Agency and dated December 29, 1993 the regional groundwater flow direction to the west of the site appears to be to the southwest. However, based upon the measured depth to water in the five wells at the site on September 29, 1992 Aegis identified a northwesterly groundwater flow direction. Based upon water level measurements collected by P&D from the five wells at the site on February 18, July 5, and October 12, 1994, February 1, and May 4, 1995 the groundwater flow direction at the site was calculated to be to the north, towards San Leandro Creek. Based upon water level measurements collected in wells MW1 through MW8 by P&D personnel on June 23 and December 19, 1995, March 28 and June 21, 1996 the groundwater flow direction was calculated to be to the northwest

The measured depth to water at or near the site on November 25, 2003 for all of the wells ranged from 38.43 to 41.70 feet. Since the previous monitoring on July 15, 2003, groundwater elevations have decreased in all of the wells by between 0.35 and 0.53 feet. The groundwater flow direction on November 25, 2003 was to the northwest with a gradient of 0.046. The calculated water level in well MW3 appears to be inconsistent with the other wells in the groundwater monitoring network.

The groundwater flow direction and gradient have remained relatively unchanged since the previous water level measurements on July 15, 2003. The groundwater monitoring data are presented in Table 1. The groundwater flow direction at the site on November 25, 2003 is shown on Figure 2.

LABORATORY RESULTS

All of the groundwater samples collected from the monitoring wells were analyzed for TPH-G using EPA Method 5030 in conjunction with Modified EPA Method 8015 and for Volatile Organic Compounds (VOCs), inleuding fuel oxygenates using EPA Method 8260 in accordance with a request from Ms. Eva Chu of the ACDEH.

The laboratory analytical results for the groundwater samples showed that TPH-G, BTEX, MTBE, and other fuel oxygenates were not detected in wells MW5, MW6, MW7, and MW8, except for 0.00084 milligrams per liter (mg/L) MTBE in well MW5, 0.00076 to 0.0014 mg/L of chloroform in wells MW6, MW7, and MW8; and 0.00078 mg/L tetrachloroethene in well MW7. TPH-G was not detected in well MW4, and was detected in wells MW1, MW2, and MW3 at concentrations of 0.140, 65, and 0.11 mg/L, respectively. In addition to well Mw5, MTBE was detected in wells MW1, MW2, MW3, and MW4 at concentrations of 0.032, 2.7, 0.33, and 8.8 mg/L, respectively. Benzene was only detected in wells MW1 and MW2 at concentrations of 0.0025 and 6.8 mg/L,

respectively. Additional compounds (toluene, ethylbenzene, xylenes, naphthalene, 1,2,4-trimethylbenzene, and 1,3,5-trimethylbenzene were also detected in well MW2.

Since the previous sampling event, TPH-G concentrations have increased in well MW1; decreased in wells MW2 and MW3; and remained unchanged (not detected) in wells MW5 through MW8. In the sample from well MW4, the reporting limit was raised for TPH-G, due to the high MTBE concentration. MTBE concentrations have increased in wells MW4 and MW5; decreased in wells MW1, MW2, and MW3; and remained unchanged (not detected) in wells MW6, MW7, and MW8. Benzene concentrations have increased in wells MW1 and MW2 and remained unchanged (not detected) in wells MW3 through MW8.

The laboratory analytical results are summarized in Table 2. Copies of the laboratory analytical report and chain of custody documentation are attached with this report.

DISCUSSION AND RECOMMENDATIONS

Based on the depth to water measurements on November 25, 2003 from all of the monitoring wells (MW1 through MW8), the groundwater flow direction at the subject site is to the northwest. The groundwater flow direction and gradient have remained relatively unchanged since the previous monitoring event. MTBE concentrations have decreased in wells MW1, MW2, and MW3; and increased in wells MW4 and MW5; remained unchanged (not detected) in wells MW6, MW7, and MW8. Benzene concentrations have increased in wells MW1 and MW2 and remained unchanged (not detected) in wells MW3 through MW8. Other BTEX and fuel oxygenates were not detected in any of the wells other than well MW2.

P&D recommends that the quarterly groundwater monitoring and sampling program be continued.

DISTRIBUTION

Copies of this report should be forwarded to Ms. Donna Drogos at the ACDEH.

LIMITATIONS

This report was prepared solely for the use of Mr. Manmohan Chopra. The content and conclusions provided by P&D in this assessment are based on information collected during our investigation, which may include, but not be limited to, visual site inspections; interviews with the site owner, regulatory agencies and other pertinent individuals; review of available public documents; subsurface exploration and our professional judgement based on said information at the time of preparation of this document. Any subsurface sample results and observations presented herein are considered to be representative of the area of investigation; however, geological conditions may vary between borings and pits and may not necessarily apply to the general site as a whole. If future subsurface or other conditions are revealed which vary from these findings, the newly revealed conditions must be evaluated and may invalidate the findings of this report.

This report is issued with the understanding that it is the responsibility of the owner, or his representative, to ensure that the information contained herein is brought to the attention of the appropriate regulatory agencies, where required by law. Additionally, it is the sole responsibility of the owner to properly dispose of any hazardous materials or hazardous wastes left onsite, in accordance with existing laws and regulations.

This report has been prepared in accordance with generally accepted practices using standards of care and diligence normally practiced by recognized consulting firms performing services of a similar nature. P&D is not responsible for the accuracy or completeness of information provided by other individuals or entities, which is used in this report. This report presents our professional judgment based upon data and findings identified in this report and interpretation of such data based upon our experience and background, and no warranty, either express or implied, is made. The conclusions presented are based upon the current regulatory climate and may require revision if future regulatory changes occur.

Should you have any questions, please do not hesitate to contact us at (510) 658-6916.

Sincerely,

P&D Environmental

Paul H. King

President

California Registered Geologist #5901

Expires: 12/31/05

Attachments: Tables 1 & 2

Site Location Map (Figure 1)

Site Plan (Figure 2) Field Parameter Forms

Laboratory Analytical Reports Chain of Custody Documentation

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TABLE 1 WELL MONITORING DATA

Well No.	Date Monitored	Top of Casing Elev. (ft.)	Depth to Water (ft.)	Water Table Elev. (ft.)
1.0.	1/10/Http://d	Diev. (It.)	Water (II.)	Diot. (it.)
MW1	11/25/03	Not Available	40.00	Not Available
	7/15/03	Not Available	39.60	Not Available
	4/16/03	Not Available	38.91	Not Available
	1/20/03	Not Available	38.21	Not Available
	2/16/99	Not Available	34.58	Not Available
	1/25/98	Not Available	33.70	Not Available
	7/14/97	Not Available	39.45	Not Available
	3/11/97	87.98+	36.90	51.08
	6/21/96		38.56	49.42
	3/28/96		37.10	50.88
	12/19/95		40.16	47.82
	6/23/95		38.54	49.44
	5/04/95	87.96++	37.65	50.33
	2/01/95		38,46	49.52
	10/12/94		42.01	45.97
	7/05/94		41.36	46.62
	2/18/94		41.02	46.96
	9/29/92		42.77	45.21
MW2	11/25/03	86.61+	38.68	47.93
	7/15/03		38.15	48.46
	4/16/03		37.50	49.11
	1/20/03		37.04	49.57
	2/16/99		33.51	53.10
	1/25/98		32.80	53.81
	7/14/97		38.46	48.15
	3/11/97		35.71	50.90
	6/21/96		37.30	49.31
	3/28/96		35.97	50.64
	12/19/95		38.80	47.81
	6/23/95		37.40	49.21
	5/04/95	86.60++	36.54	50.07
	2/01/95		37.27	49.34
	10/12/94		40.77	45,84
	7/05/94		40.13	46.48
	2/18/94		39.81	46.80
	9/29/92		41.55	45.06

NOTES:

ft. = Feet.

^{+ =} Indicates survey data provided by Kier & Wright dated June 26, 1995.

^{++ =} Indicates survey data provided by Aegis Environmental, Inc.

Well No.	Date Monitored	Top of Casing	Depth to Water (ft.)	Water Table Elev. (ft.)
110.	Monnoten	Elev. (ft.)	water (it.)	Elev. (II.)
MW3	11/25/03	87.48+	41.70	45.78
	7/15/03		41.34	46.14
	4/16/03		40.60	46.88
	1/20/03		39.81	47.67
	2/16/99		34.91	52,57
	1/25/98		33.91	53,57
	7/14/97		40.61	46.87
	3/11/97		38.71	48.77
	6/21/96		40.61	46.87
	3/28/96		38.75	48.73
	12/19/95		42.20	45.28
	6/23/95		40.65	46.83
	5/04/95	87.50++	39.61	47.87
	2/01/95		40.13	47.35
	10/12/94		43.92	43.56
	7/05/94		43.32	44.16
	2/18/94		43.09	44.39
	9/29/92		44.60	42.88*
MW4	11/25/03	86.21+	38.43	47.78
	7/15/03		38.04	48.17
	4/16/03		37.32	48.89
	1/20/03		36.70	49.51
	2/16/99		33.43	52.78
	1/25/98		32.96	53.25
	7/14/97		38,10	48.11
	3/11/97		33.24	52.97
	6/21/96		37.12	49.09
	3/28/96		35,00	51.21
	12/19/95		38.45	47.76
	6/23/95		37.40	48,81
	5/04/95	86.20++	36.33	49.88
	2/01/95		36.96	49.25
	10/12/94		40.48	45.7 3
	7/05/94		39.69	46,52
	2/18/94		39.36	46.85
	9/29/92		44.29	41.92

NOTES:

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^{++ =} Indicates survey data provided by Aegis Environmental, Inc.

^{** =} Indicates depth to water measurements prior to groundwater monitoring well development.

Well No.	Date Monitored	Top of Casing Elev. (ft.)	Depth to Water (ft.)	Water Table Elev. (ft.)
			·· ==== (==,	
MW5	11/25/03	89.10+	41.41	47.69
	7/15/03		41.06	48.04
	4/16/03		39.92	49.18
	1/20/03		39.50	49.60
	2/16/99		35.08	54.02
	1/25/98		34.08	55,02
	7/14/97		41.20	47.90
	3/11/97		38.02	51,08
	6/21/96		40.03	49.07
	3/28/96		38.30	50.80
	12/19/95		41.79	47.31
	6/23/95		39.87	49.23
	5/04/95	89.06++	38.94	50,16
	2/01/95		39,94	49.16
	10/12/94		43,81	45.29
	7/05/94		43.08	46.02
	2/18/94		42,88	46.22
	9/29/92		44.53	44.57
MW6	11/25/03	84.02+	38.97	45.05
	7/15/03		38.61	45.41
	4/16/03		38.00	46.02
	1/20/03		37.21	46.81
	2/16/99		32.82	51.20
	1/25/98		31.64	52.38
	7/14/97		39.04	44.98
	3/11/97		36.32	47.70
	6/21/96		38.00	46.02
	3/28/96		36.18	47.84
	12/19/95		39.25	44.77
	6/23/95		38.17	44.77 45.85
	6/21/95**			
	U/Z1/73 ***		38.11	45.91

NOTES:

ft. = Feet.

^{+ =} Indicates survey data provided by Kier & Wright dated June 26, 1995.

^{++ =} Indicates survey data provided by Aegis Environmental, Inc.

^{** =} Indicates depth to water measurements prior to groundwater monitoring well development.

Well	Date	Top of Casing	Depth to	Water Table
No.	Monitored	Elev. (ft.)	Water (ft.)	Elev. (ft.)
MW7	11/25/03	87.11+	41.68	45.43
	7/15/03		41.30	45.81
	4/16/03		40.63	46,48
	1/20/03		39. <i>77</i>	47.34
	2/16/99		34.59	52.52
	1/25/98		33.47	53.64
	7 /14/97		41.97	45.14
	3/11/97		38.96	48.15
	6/21/96		40.80	46.31
	3/28/96		38.94	48.17
	12/19/95		42.26	44.85
	6/23/95		41.00	46.11
	6/21/95**	•	40.30	46.81
MW8	11/25/03	89.70+	40.92	48.78
	7/15/03		40.50	49.20
	4/16/03		39.52	50.18
	1/20/03		38.94	50.76
	2/16/99		33,92	55.78
	1/25/98		32.73	56.97
	7/14/97		39.98	49.72
	3/11/97		36.74	52.96
	6/21/96		38.69	51.01
	3/28/96		36.98	52.72
	12/19/95		40,35	49.35
	6/23/95		38.36	51.34
	6/21/95**		38.20	51.50

NOTES:

ft = Feet

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^{** =} Indicates depth to water measurements prior to groundwater monitoring well development.

TABLE 2 GROUNDWATER LABORATORY ANALYTICAL RESULTS (MW1)

Date	TPH-G	MTBE	Benzene	Toluene	Ethyl- benzene	Total Xylenes	Other VOCs by EPA 8260
11/25/03	0.140	0.032	0.0025	ND<0.001	ND<0.001	ND<0.001	ND
07/15/03	0.060	0.053	ND<0,001	ND<0.001	ND<0.001	ND<0.001	ND, except t-Butyl Alcohol (TBA) = 0.012
04/17/03	0.052	0.056	0.0011	ND<0.001	ND<0.001	ND<0.001	ND, except t-Butyl Alcohol (TBA) = 0.013
01/20/03	0.17	0.085	ND<0.005	ND<0.005	ND<0.005	ND<0.005	ND
02/17/99	0.97	0.29	0.067	0.12	0.0093	0.058	
01/25/98	0.30	ND<0.014	0.021	0.00073	0.0076	0.0010	
07/14/97	0.20	0.035	0.020	0.0055	0.0012	0.0023	
03/11/97	0.60	0.014	0.053	0.00095	0.003	0.0015	
06/21/96	1.4	0.019	0.30	0.0087	0.033	0.0098	
03/28/96	1.3	0.022	0.32	0.0023	0.034	0.0046	
12/19/95	0.50	0.0081	0.087	0.0015	0.011	0.0035	
06/23/95	Not	Sampled					
05/4/95	2,4		0.67	0.0028	0.076	0.0060	**
02/01/95	4.6		1.8	0.0099	0.23	0.030	
10/12/94	2.5		0.82	0.0039	0.10	0.020	
07/05/94	3.0		1.3	0.0038	0.035	0.0025	
09/29/92	3.1	•	0.16	ND	ND	0.0060	

NOTES:

TPH-G = Total Petroleum Hydrocarbons as Gasoline.

MTBE = Methyl Tert Butyl Ether.

VOCs = Volatile Organic Compounds

ND = Not Detected.

-- = Sample not analyzed for this compound during this sampling event.

TABLE 1 WELL MONITORING DATA

Well Date Top of Casing Depth to No. Monitored Elev. (ft.) Water (ft.) MW1 11/25/03 Not Available 40.00 7/15/03 Not Available 39.60	Elev. (ft.) Not Available Not Available Not Available Not Available
	Not Available Not Available
7/15/03 Not Available 20.40	Not Available
// 1.5/05 INDUAVALIABLE 57,00	
4/16/03 Not Available 38,91	Not Available
1/20/03 Not Available 38.21	
2/16/99 Not Available 34.58	Not Available
1/25/98 Not Available 33.70	Not Available
7/14/97 Not Available 39.45	Not Available
3/11/97 87.98+ 36,90	51.08
6/21/96 38.56	49.42
3/28/96 37.10	50.88
12/19/95 40.16	47.82
6/23/95 38.54	49.44
5/04/95 87.96++ 37.65	50.33
2/01/95 38.46	49.52
10/12/94 42.01	45.97
7/05/94 41.36	46.62
2/18/94 41.02	46.96
9/29/92 42.77	45.21
MW2 11/25/03 86.61+ 38.68	47.93
7/15/03 38.15	48,46
4/16/03 37.50	49.11
1/20/03 37.04	49.57
2/16/99 33.51	53.10
1/25/98 32.80	53.81
7/14/97 38.46	48.15
3/11/97 35.71	50.90
6/21/96 37.30	49,31
3/28/96 35,97	50.64
12/19/95 38,80	47.81
6/23/95 37.40	49.21
5/04/95 86.60++ 36,54	50.07
2/01/95 37.27	49.34
10/12/94 40.77	45.84
7/05/94 40.13	46.48
2/18/94 39.81	46.80
9/29/92 41.55	45.06

NOTES:

ft. = Feet

^{+ =} Indicates survey data provided by Kier & Wright dated June 26, 1995.

^{++ =} Indicates survey data provided by Aegis Environmental, Inc.

Well No.	Date Monitored	Top of Casing Elev. (ft.)	Depth to Water (ft.)	Water Table Elev. (ft.)
MW3	11/25/03	87.48+	41.70	45.78
	7/15/03		41.34	46.14
	4/16/03		40.60	46.88
	1/20/03		39.81	47.67
	2/16/99		34.91	52.57
	1/25/98		33.91	53,57
	7/14/97		40.61	46.87
	3/11/97		38.71	48.77
	6/21/96		40.61	46.87
	3/28/96		38.75	48.73
	12/19/95		42.20	45.28
	6/23/95		40.65	46.83
	5/04/95	87.50++	39.61	47.87
	2/01/95		40.13	47.35
	10/12/94		43.92	4 3. 5 6
	7/05/94		43.32	44.16
	2/18/94		43,09	44.39
	9/29/92		44.60	42.88*
MW4	11/25/03	86.21+	38.43	47.78
	7/15/03		38.04	48.17
	4/16/03		37.32	48.89
	1/20/03		36.70	49,51
	2/16/99		33.43	52.78
	1/25/98		32.96	53,25
	7/14/97		38.10	48.11
	3/11/97		33.24	52.97
	6/21/96		37.12	49.09
	3/28/96		35.00	51.21
	12/19/95		38.45	47.76
	6/23/95		37.40	48.81
	5/04/95	86.20++	36.33	49.88
	2/01/95		36,96	49.25
	10/12/94		40.48	45.73
	7/05/94		39.69	46.52
	2/18/94		39.36	46.85
	9/29/92		44.29	41.92

NOTES:

ft. = Feet.

^{+ =} Indicates survey data provided by Kier & Wright dated June 26, 1995.

^{++ =} Indicates survey data provided by Aegis Environmental, Inc.

^{** =} Indicates depth to water measurements prior to groundwater monitoring well development.

TABLE 1 (Continued)
WELL MONITORING DATA

Well No.	Date Monitored	Top of Casing Elev. (ft.)	Depth to Water (ft.)	Water Table Elev. (ft.)
MW5	11/25/03	89.10+	41.41	47.69
	7/15/03		41.06	48.04
	4/16/03		39.92	49.18
	1/20/03		39.50	49.60
	2/16/99		35.08	54.02
	1/25/98		34.08	55.02
	7/14/97		41.20	47.90
	3/11/97		38.02	51.08
	6/21/96		40.03	49.07
	3/28/96		38,30	50.80
	12/19/95		41.79	47.31
	6/23/95		39.87	49.23
	5/04/95	89.06++	38.94	50,16
	2/01/95	0 7.00 · ·	39.94	49.16
	10/12/94		43.81	45.29
	7/05/94		43.08	46.02
	2/18/94		42.88	46.22
	9/29/92		44.53	44.57
	7/27/72		44,55	44.07
MW6	11/25/03	84.02+	38.97	45,05
	7/15/03		38.61	45.41
	4/16/03		38.00	46.02
	1/20/03		37.21	46.81
	2/16/99		32.82	51.20
	1/25/98		31.64	52.38
	7/14/97		39.04	44,98
	3/11/97		36.32	47.70
	6/21/96		38.00	46.02
	3/28/96		36.18	47.84
	12/19/95		39.25	44.77
	6/23/95		38.17	45.85
	6/21/95**		38.11	45.91
	V. = 2, 2 V		20.11	

NOTES:

ft. = Feet.

^{+ =} Indicates survey data provided by Kier & Wright dated June 26, 1995.

^{+++ =} Indicates survey data provided by Aegis Environmental, Inc.

^{** =} Indicates depth to water measurements prior to groundwater monitoring well development.

Well No.	Date Monitored	Top of Casing Elev. (ft.)	Depth to Water (ft.)	Water Table Elev. (ft.)
MW7	11/25/03	87.11+	41.68	45.43
	7/15/03		41.30	45.81
	4/16/03		40.63	46.48
	1/20/03		39.77	47.34
	2/16/99		34.59	52.52
	1/25/98		33.47	53.64
	7/14/97		41.97	45.14
	3/11/97		38.96	48.15
	6/21/96		40.80	46.31
	3/28/96		38.94	48.17
	12/19/95		42.26	44.85
	6/23/95		41.00	46.11
	6/21/95**		40.30	46.81
MW8	11/25/03	89.70+	40.92	48.78
	7/15/03		40.50	49.20
	4/16/03		39.52	50.18
	1/20/03		38.94	50.76
	2/16/99		33.92	55,78
	1/25/98		32.73	56.97
	7/14/97		39.98	49.72
	3/11/97		36.74	52.96
	6/21/96		38.69	51.01
	3/28/96		36.98	52.72
	12/19/95		40.35	49.35
	6/23/95		38.36	51.34
	6/21/95**		38.20	51.50

NOTES:

ft. = Feet.

^{+ =} Indicates survey data provided by Kier & Wright dated June 26, 1995.

^{** =} Indicates depth to water measurements prior to groundwater monitoring well development.

TABLE 2 GROUNDWATER LABORATORY ANALYTICAL RESULTS (MW1)

Date	TPH-G	MTBE	Benzene	Toluene	Ethyl- benzene	Total Xylenes	Other VOCs by EPA 8260
11/25/03	0.140	0.032	0.0025	ND<0.001	ND<0.001	ND<0.001	ND
07/15/03	0.060	0.053	ND<0.001	ND<0.001	ND<0.001	ND<0.001	ND, except t-Butyl Alcohol (TBA) = 0.012
04/17/03	0.052	0.056	0.0011	ND<0.001	ND<0.001	ND<0.001	ND, except t-Butyl Alcohol (TBA) = 0.013
01/20/03	0.17	0.085	ND<0.005	ND<0.005	ND<0.005	ND<0.005	ND
02/17/99	0.97	0.29	0.067	0.12	0.0093	0.058	
01/25/98	0.30	ND<0.014	0.021	0.00073	0.0076	0.0010	•
07/14/97	0.20	0.035	0.020	0.0055	0.0012	0.0023	
03/11/97	0.60	0.014	0.053	0.00095	0.003	0.0015	
06/21/96	1.4	0.019	0.30	0.0087	0.033	0.0098	
03/28/96	1.3	0.022	0.32	0.0023	0.034	0.0046	
12/19/95	0.50	0.0081	0.087	0.0015	0.011	0.0035	
06/23/95	Not	Sampled					
05/4/95	2.4		0.67	0.0028	0.076	0.0060	
02/01/95	4.6		1.8	0.0099	0.23	0.030	-
10/12/94	2.5		0.82	0.0039	0.10	0.020	
07/05/94	3.0		1.3	0.0038	0.035	0.0025	
09/29/92	3.1		0.16	ND	ND	0.0060	•••

NOTES:

TPH-G = Total Petroleum Hydrocarbons as Gasoline.

MTBE = Methyl Tert Butyl Ether.

VOCs = Volatile Organic Compounds

ND = Not Detected.

-- = Sample not analyzed for this compound during this sampling event.

TABLE 2 (Continued) GROUNDWATER LABORATORY ANALYTICAL RESULTS (MW2)

Date	TPH-G	МТВЕ	Benzene	Toluene	Ethyl- benzene	Total Xylenes	Other VOCs by EPA 8260
11/25/03	65	2.7	6.8	8,8	2.9	16	ND, except Naphtalene = 0.54 1,2,4-Trimethylbenzene = 1.8 1,3,5-Trimethylbenzene = 0.42
07/15/03	78	4.1	3.3	4.4	1.8	9.3	ND, except Naphthalene = 0.29 1,2,4-Trimethylbenzene = 1.3 1,3,5-Trimethylbenzene = 0.32
04/17/03	57	5.6	3.4	5.1	2.8	10	n-Propyl benzene = 0.15 ND, except Naphthanlene = 0.43 1,2,4-Trimethylbenzene = 2.2 n-proylbenzene = 0.26 1,3,5-Trimethylbenzene = 0.55
01/20/03	48	3.8	2.9	3.0	2.0	11	ND, except Naphthanlene = 0.35 1,2,4-Trimethylbenzene = 1.4 1,3,5-Trimethylbenzene = 0.32 Isopropylbenzene = 0.069 n-Propyl benzene = 0.16
02/17/99	7.3	0.29	0.067	0.12	0.0093	0.058	
01/25/98	24	2.7	2.7	4.9	0.70	4.0	
07/14/97	43	1.6	6.2	8.9	1.5	7.4	
03/11/97	28	0.71	4.0	4.5	0.99	4.3	
06/21/96	49	0.53	6.6	6.3	1.4	6.2	
03/28/96	38	0.45	5.8	4.7	1.1	5.1	
12/19/95	25	0.45	5.2	3.8	0.86	3.8	
06/23/95	Not	Sampled					
05/4/95	63		10	11	1.6	8.8	
02/01/95	45		7.0	5.1	1.2	6.1	
10/12/94	24		4.4	2.8	0.73	3.5	
07/05/94	46		9.1	7.0	1.4	7.3	
09/29/92	20	-	4.6	3.8	0.26	3.3	

NOTES:

TPH-G = Total Petroleum Hydrocarbons as Gasoline.

MTBE = Methyl Tert Butyl Ether.

VOCs = Volatile Organic Compounds

ND = Not Detected.

--= Sample not analyzed for this compound during this sampling event.

--= Sample not analyzed for this compound during this sampling order.

Results are reported in milligrams per kilograms (mg/L), unless otherwise specified.

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TABLE 2 (Continued) GROUNDWATER LABORATORY ANALYTICAL RESULTS (MW3)

Date	TPH-G	МТВЕ	Benzene	Toluene	Ethyl- benzene	Total Xylenes	Other VOCs by EPA 8260
11/25/03	0.11	0.33	ND<0.005	ND<0.005	ND<0.005	ND<0.005	ND
07/15/03	0.16	0.66	ND<0.0012	ND<0.0012	ND<0.0012	ND<0.0012	ND
							. —
04/17/03	0.18	0.34	ND	ND	ND	ND	ND
01/20/03	0.12	0.25	ND<0.005	ND<0.005	ND<0.005	0.0052	ND
02/17/99	ND	0.29	0.067	0.12	0.0093	0,058	
01/25/98	0.49	0.71	0.0079	0,0061	0.0053	0.029	
07/14/97	0.40	0.11	0.00093	0.010	0.0013	0.00068	
03/11/97	1.1	0.68	0.053	0.013	0.063	0.017	
06/21/96	1.3	0.3	0.094	0.0021	0.039	0.002	
03/28/96	4.6	1.1	1.4	0.012	0.17	0.020	
12/19/95	0.95	0.12	0.16	0.0023	0.015	0.0016	
06/23/95	Not	Sampled					
05/4/95	7.2		3.1	0.038	0.20	0.062	
02/01/95	11		4.2	0.031	0.33	0.29	**
10/12/94	1.7		0.39	0.00090	0.018	0.0057	
07/05/94	3.6		1.6	0.0083	0.076	0.047	
09/29/92	Not	Sampled					

NOTES:

TPH-G = Total Petroleum Hydrocarbons as Gasoline.

MTBE = Methyl Tert Butyl Ether.

VOCs = Volatile Organic Compounds

ND = Not Detected.

-- = Sample not analyzed for this compound during this sampling event.

TABLE 2 (Continued) GROUNDWATER LABORATORY ANALYTICAL RESULTS (MW2)

Date	TPH-G	MTBE	Benzene	Toluene	Ethyl- benzene	Total Xylenes	Other VOCs by EPA 8260
11/25/03	65	2.7	6.8	8.8	2.9	16	ND, except Naphtalene = 0.54 1,2,4-Trimethylbenzene = 1.8 1,3,5-Trimethylbenzene = 0.42
07/15/03	78	4.1	3.3	4.4	1.8	9.3	ND, except Naphthalene = 0.29 1,2,4-Trimethylbenzene = 1.3 1,3,5-Trimethylbenzene = 0.32 n-Propyl benzene = 0.15
04/17/03	57	5.6	3.4	5.1	2.8	10	ND, except Naphthanlene = 0.43 1,2,4-Trimethylbenzene = 2.2 n-proylbenzene = 0.26 1,3,5-Trimethylbenzene = 0.55
01/20/03	48	3.8	2.9	3.0	2.0	11	ND, except Naphthanlene = 0.35 1,2,4-Trimethylbenzene = 1.4 1,3,5-Trimethylbenzene = 0.32 Isopropylbenzene = 0.069 n-Propyl benzene = 0.16
02/17/99	7.3	0.29	0.067	0.12	0.0093	0.058	
01/25/98	24	2.7	2.7	4.9	0.70	4.0	
07/14/97	43	1.6	6.2	8.9	1.5	7.4	
03/11/97	28	0.71	4.0	4.5	0.99	4.3	*
06/21/96	49	0.53	6.6	6.3	1.4	6.2	
03/28/96	38	0.45	5.8	4.7	1.1	5.1	
12/19/95	25	0.45	5.2	3.8	0.86	3.8	-
06/23/95	Not	Sampled					
05/4/95	63		10	11	1.6	8.8	-
02/01/95	45		7.0	5.1	1.2	6.1	
10/12/94	24		4.4	2.8	0.73	3.5	
07/05/94	46		9.1	7.0	1.4	7.3	
09/29/92	20		4.6	3.8	0.26	3.3	

NOTES:

TPH-G = Total Petroleum Hydrocarbons as Gasoline.

MTBE = Methyl Tert Butyl Ether.

VOCs = Volatile Organic Compounds

ND = Not Detected.

-- = Sample not analyzed for this compound during this sampling event.

TABLE 2 (Continued) GROUNDWATER LABORATORY ANALYTICAL RESULTS (MW5)

Date	ТРН-G	MTBE	Benzene	Toluene	Ethyl- benzene	Total Xylenes	Other VOCs by EPA 8260
11/25/03	ND	0.00084	ND<0.001	ND<0.001	ND<0.001	ND<0.001	ND
07/15/03	ND	0.0014	ND	ND	ND	ND	ND
04/17/03	ND	ND	ND	ND	ND	ND	ND
01/20/03	ND	ND	ND	ND	ND	ND	ND
02/17/99	0.17	ND	ND	0.00074	ND	ND	
01/25/98	ND	ND	ND	ND	ND	ND	
07/14/97	ND	ND	ND	ND	ND	ND	
03/11/97	ND	ND	ND	ND	ND	0.00077	
06/21/96	ND	ND	ND	ND	ND	ND	+-
03/28/96	ND	ND	ND	ND	ND	ND	
12/19/95	ND	ND	ND	ND	ND	ND	
06/23/95	Not	Sampled					
05/4/95	ND		ND	ND	ND	ND	
02/01/95	ND		ND	ND	ND	ND	
10/12/94	ND		ND	ND	ND	ND	
07/05/94	ND		ND	ND	ND	0.0010	
09/29/92	0.06		10	0.0071	ND	0.0069	

NOTES:

TPH-G = Total Petroleum Hydrocarbons as Gasoline.

MTBE = Methyl Tert Butyl Ether.

VOCs = Volatile Organic Compounds

ND = Not Detected.

-- = Sample not analyzed for this compound during this sampling event.

TABLE 2 (Continued) GROUNDWATER LABORATORY ANALYTICAL RESULTS (MW6)

Date	TPH-G	MTBE	Benzene	Toluene	Ethyl- benzene	Total Xylenes	Other VOCs by EPA 8260
11/25/03	ND<0.05	ND<0.001	ND<0.001	ND<0.001	ND<0.001	ND<0.001	ND, except Chloroform = 0.00089
07/15/03	ND	ND	ND	ND	ND	ND	ND, except Chloroform = 0.00084 1,2-Dibromo- 3-chloropropane = 0.00066 Tetrachloroethene = 0.00067
04/17/03	ND	ND	ND	ND	ND	ND	ND, except Chloroform = 0.0012
01/20/03	ND	0.0012	ND	ND	ND	ND	ND, except Chloroform = 0.0011
02/17/99	ND	ND	ND	ND	ND	ND	
01/25/98	ND	ND	ND	ND	ND	ND	
07/14/97	ND	0.019	ND	ND	ND	ND	-
03/11/97	ND	ND	ND	ND	ND	ND	_
06/21/96	ND	ND	ND	ND	ND	ND	
03/28/96	ND	ND	ND	ND	ND	ND	
12/19/95	ND	0.01	ND	ND	ND	ND	
06/23/95	ND	0.003	ND	ND	ND	ND	

NOTES:

TPH-G = Total Petroleum Hydrocarbons as Gasoline.

MTBE = Methyl Tert Butyl Ether.

VOCs = Volatile Organic Compounds

ND = Not Detected.

-- = Sample not analyzed for this compound during this sampling event.

TABLE 2 (Continued) GROUNDWATER LABORATORY ANALYTICAL RESULTS (MW7)

Date	TPH-G	MTBE	Benzene	Toluene	Ethyl- benzene	Total Xylenes	Other VOCs by EPA 8260
11/25/03	ND<0.05	ND<0.001	ND<0.001	ND<0.001	ND<0.001	ND<0.001	ND, except Chloroform = 0.00076
07/15/03	ND	ND	ND	ND	ND	ND	Tetrachloroethene = 0.00078 ND, except Chloroform = 0.00061 1,2-Dibromo- 3-chloropropane = 0.00064
04/17/03	ND	ND	ND	ND	ND	ND	Tetrachloroethene = 0.0012 ND, except Chloroform = 0.00075 Tetrachloroethene = 0.0012
01/20/03	ND	ND	ND	ND	ND	ND	ND, except Chloroform = 0.00056
02/17/99	ND	ND	ND	ND	ND	ND	==
01/25/98	ND	ND	ND	ND	ND	ND	
07/14/97	ND	ND	ND	ND	ND	ND	
03/11/97	ND	ND	ND	ND	ND	ND	
06/21/96	ND	ND	ND	ND	ND	ND	
03/28/96	ND	ND	ND	ND	ND	ND	## .
12/19/95	ND	ND	ND	ND	ND	ND	
06/23/95	ND	ND	ND	ND	ND	ND	

NOTES:

TPH-G = Total Petroleum Hydrocarbons as Gasoline.

MTBE = Methyl Tert Butyl Ether.

VOCs = Volatile Organic Compounds

ND = Not Detected.

-- = Sample not analyzed for this compound during this sampling event.

TABLE 2 (Continued) GROUNDWATER LABORATORY ANALYTICAL RESULTS (MW8)

Date	TPH-G	MTBE	Benzene	Toluene	Ethyl- benzene	Total Xylenes	Other VOCs by EPA 8260
11/25/03	ND<0.05	ND<0.001	ND<0,001	ND<0.001	ND<0.001	ND<0.001	ND, except Chloroform = 0.0014
07/15/03	ND	ND	ND	ND	ND	0.00066	ND, except Chloroform = 0.0014 1,2-Dibromo- 3-chloropropane = 0.00052
04/17/03	ND	ND	ND	ND	ND	ND	ND, except Chloroform = 0.0018
01/20/03	ND	ND	ND	ND	ND	ND	ND, except Chloroform = 0.0013
02/17/99	ND	ND	ND	ND	ND	ND	
01/25/98	ND	ND	ND	ND	ND	ND	
07/14/97	ND	ND	ND	ND	ND	ND	
03/11/97	ND	ND	ND	ND	ND	ND	
06/21/96	ND	ND	ND	ND	ND	ND	-
03/28/96	ND	ND	ND	ND	ND	ND	
12/19/95	ND	ND	ND	ND	ND	ND	
06/23/95	ND	ND	ND	ND	ND	ND	

NOTES:

TPH-G = Total Petroleum Hydrocarbons as Gasoline.

MTBE = Methyl Tert Butyl Ether.

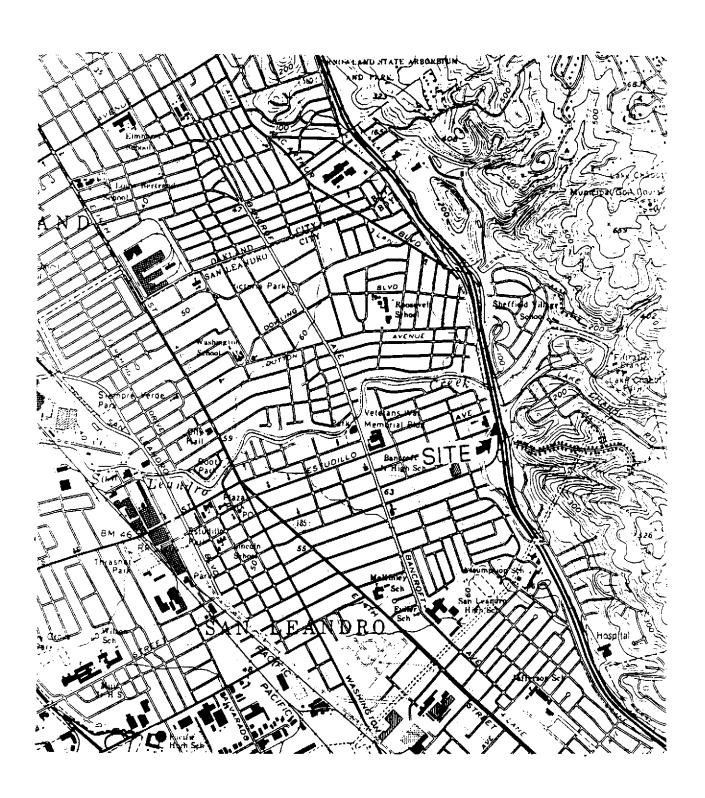
VOCs = Volatile Organic Compounds

ND = Not Detected.

-- = Sample not analyzed for this compound during this sampling event.

P & D Environmental

4020 Panama Court Oakland, CA 94611 Telephone (510) 658-6916



Base Map from: U.S. Geological Survey San Leandro, Calif. 7.5 Minute Quadrangle Photorevised 1980

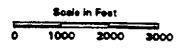
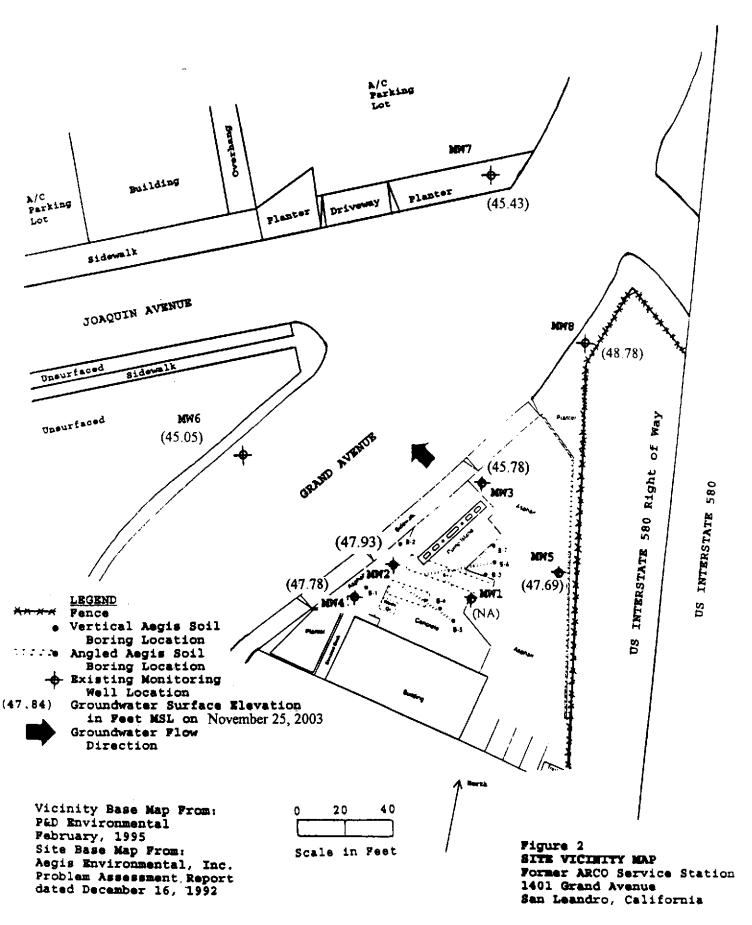


Figure 1
SITE LOCATION MAP
Former ARCO Service Station
1401 Grand Avenue
San Leandro, California

P & D ENVIRONMENTAL

4020 Panama Court Oakland, CA 94611 Telephone (510) 658-6916



sice Name Former Haber Oil	Well No. MW
Job No. 0655	Date 1/25/03
TOC to Water (ft.) 40.00	Sheen NONE
Well Depth (ft.) 52.7	Free Product Thickness
Well Diameter <u>Uin</u>	Sample Collection Method
Gal./Casing Vol. 8.2	teflor bailer
Z=24.6 TIME GAL, PURGED PH	TEMPERATURE ELECTRICAL (45/Cm) X/C
2'.00 3 7.84	62.0 0.64
	62.9 0.62
2:04 8 7.60 2:08 12 7.60	63.0 0.65
2:12 15 7.64	62.8 0.65
2:17 10 8:038	62.4 0.63
2:22 25 7.74	62.8 0.62
7/27	600
- Santing Inc.	
	
	
	· · · · · · · · · · · · · · · · · · ·
<u> </u>	
	1.1
NOTES: & pH value did not Very Slight PHC odor on	Stabilize at this time
Very Slight PHC oder on	bural water
PURGE10.92	

Site Name Former Hober Cil	Well No. MWZ
JOD NO. 6655	Date 11/25/03
TOC to Water (ft.) 38.68	Sheen NONE
Well Depth (ft.) 52.4	Free Product Thickness
Well Diameter + in .	Sample Collection Method
Gal./Casing Vol. 8,9	Teflon bailer
$\mathcal{E} = 26.7$ TIME GAL PURGED PH	TEMPERATURE CONDUCTIVITY
3:59 5 7.27	63.5 C.73
4.03 10 7.24	65.0 C,72
4:07 15 7.34	64,6 0,74
4:12 20 7.39	64.3 6.71
4'15 23 7.41	65.2 0.72
4.18 27 7.38	69.1 0.74
4:25 Sampling Ame	
<u> </u>	
	<u> </u>
NOTES: Sia OHC In I	
Strong PTC odor, b.	ut no sheen on
purge water.	
PURGE10.92	

Site Name _	Former Holbe	_0i\	Well No	mw3
Job No			Date	125/03
	r (ft.) 41,70	_	Sheen	NONE
Well Depth	(ft.) <u>55.3</u>	_	Free Produc	t Thickness
Well Diamec	er 4in	- 	· · · · · · · · · · · · · · · · · · ·	ection Method
Gal./Casing	vo1. 8,8	_	Teflo	n bailer
TIME	E=ZEM GAL. PURGED	_11	TEMPERATURE	ELECTRICAL (4 5/cm) X/CO
TIME : 05	SAL. PORGED	퍼 7.83	623	0.64
1'.(0	10	756	62.0	0.63
1:16	15	7.55	62,3	0.62
1,21	20	755	62.7	0,63
1.24	23	755	629	0.63
1:27	27	7.54	62.8	0,63
1:35	Samoling	time		
				
			·	
 .				
			· · · · · · · · · · · · · · · · · · ·	
				
				
NOTES: MØ	lone we		1	
7	tight PHC	odor, b	int no she	en on
P	nige water	<u>~; </u>		

sice Name former Haber Oil	Well No. MWY
Job No. 0055	Date 11/25/03
TOC to Water (ft.) 38,43	Sheen NONE
Well Depth $(ft.) = 53.3$	Free Product Thickness 🗸
Well Diameter <u>Hin</u>	Sample Collection Method
Gal./Casing Vol. 9.6	Jetlon bailer
£=,28,8	(OC) ELECTRICATED S/C Y/CO
TIME GAL PURGED DH	1
3:07 <u>5</u> <u>7.68</u> 3:11 10 7.63	
$\frac{7.59}{3.15}$ $\frac{7.59}{15}$	$\frac{63.9}{63.9}$ $\frac{6.87}{0.85}$
$\frac{319}{3.19}$ $\frac{7.60}{7.60}$	64.2 0.86
3: 23 <u>25</u> 7.57	64.1 0.83
7:57	
3:35 <u>29</u> 7.56	<u>4.0</u> <u>0.89</u>
- me	
	
NOTES: No DIA	1
NO 111 Codor 0	r Shan on
purge Water.	
PURGE10.92	

Site Name Former HowerCil	Well No. MW5
Job No. <u>6055</u>	Date 11/25/03
TOC to Water (ft.) 4(.4)	Sheen None
Well Depth (ft.) 54.7	Free Product Thickness
Well Diameter 4in	Sample Collection Method
Gal./Casing Vol. 4,6	Joffon bailer
∑= 25.8 TIME GAL, PURGED DH	TEMPERATURE CONDUCTIVITY (CO. * (CO.
12'13 2 7.83	647 Oil8
12:06 5 7.60	64,3 0,71
12:11 10 7.58	64.8 0.71
12:16 15 7.61	64,2 0,67
121,21 20 7,74	63,5 0,64
12:27 76 7.70	64.4 0.64
12:35 Sembline Hine	
	· · · · · · · · · · · · · · · · · · ·
HOTES: No AHC whom or	Sheen on Phrge
nicy er.	or or pringer
<u> </u>	

sice Name Former taker Oil	Well No. MW6
JOB NO. 0055	Date 11/25/03
TOC to Water (ft.) 38.97	Sheen None
Well Depth (ft.) 4911	Free Product Thickness ${\mathscr D}$
Well Diameter Zin	Sample Collection Method
Gal./Casing Vol. 1.7	Teflon Valler
S=\$5.\ TIME GAL, PURGED DH	TEMPERATURE OF ELECTRICAL (45/cm) × (0
10.53 (811)	67,3 0,64
16:54 2 7.84	64.0 0.67
10:55 3 7.83	64.4 0.67
10.56 4 7.83	64.6 0.67
16:57 5 7.81	64.5 066
10:57 6 780	645 0166
11:25 Samping the	
	· · · · · · · · · · · · · · · · · · ·
<u> </u>	
OTES: No PH Code	so or Sheen on whistic box
urge nact.	librate in christic hox
URGE10.92 Delaw TO.C.	

Sice Name Fringer Wiber Oil	Well No. MW7
Job No. 8055	Date 11/25/03
TOC to Water (ft.) 41,68	Sheen NONE
Well Depth (ft.) 49.8	Free Product Thickness
Well Diameter Zin	Sample Collection Method
Gal./Casing Vol. 1.3	Teflon bailer
TIME GAL PURGED DH	TEMPERATURE ELECTRICAL S/Cm) × Conductivity
10:19 0.5 7.90	581 0.56
10:161 1.0 7.83	614 0,59
10.20 1.5 7.80	62.2 0.61
10:21 2.2 7.83	62,5 0,61
<u>10:22</u> <u>3.0</u> <u>7.88</u>	62.5 0.63
10:22 4.C 7.85	63.162.8 062
10:201 Sounding the	
	<u> </u>
	
NOTES:	
MOTES: No PHIC ador or	Shien on
- Prige water	
PURGE10.92	

	0
sice Name former take oil	Well No. MWS
Job No. <u>CC55</u>	Date 11/25/03
TOC to Water (ft.) 40.92	Sheen NONE
Well Depth (ft.) 48,0	Free Product Thickness 💯
Well Diameter 211,	Sample Collection Method
Gal./Casing Vol	Tetlon bailer
2=3.3	TEMPERATURE (OF) ELECTRICAL (45/4) × 100
TIME GAL PURGED DH	50,100,111,1
4.36 0.5 8.12	515 0.54
$\frac{9.36}{9.37}$ 1.5 $\frac{8.00}{798}$	<u>58.7</u> <u>O.E.\</u>
	60.6 0.64
	61.1 0.64
$\frac{9.38}{2.8}$ $\frac{7.99}{2.99}$	61.3 0.64
9.39 3.5 8.00	61.3 0.64
9:45 Sampling In	·
	·
<u> </u>	
	· · · · · · · · · · · · · · · · · · ·
NOTES: Nã PHC odar	Ac Cla
on organ hader.	or shoen
PURGE10.92	

	McCampbell	Analytical	Inc.
_			

110 2nd Avenuc South, #D7, Pacheco, CA 94553-5560 Telephone: 925-798-1620 Fax: 925-798-1622 http://www.mccampbell.com E-mail: main@mccampbell.com

P & D Environmental	· · · · · · · · · · · · · · · · · · ·	Date Sampled: 11/25/03
4020 Panama Court	Oil	Date Received: 11/26/03
O-11 J. CA. 04611, 4021	Client Contact: Paul King	Date Extracted: 12/02/03-12/03/03
Oakland, CA 94611-4931	Client P.O.:	Date Analyzed: 12/02/03-12/03/03

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

raction method: S	W5030B		Analytical methods: 8015Cm	Work Order:	031138
Lab ID	Client ID	Matrix	TPH(g)	DF	% S
001A	MW I	w	140,a	1	113
002A	MW 2	w	65,000,a	50	103
003A	MW 3	w	110,m	1	113
004A	MW 4	w	ND<1000.j	10	101
005A	MW 5	w	ND	1	99.
006A	MW 6	w	ND	1	98.
007A	MW 7	w	ND	1	10
008A	MW 8	w	ND	1	10
	.				
!					
· · · · · · · · · · · · · · · · · · ·					
1					
				<u>:</u>	<u> </u>
	Limit for DF =1; not detected at or	W	50		g/L
	e reporting limit	S	NA	N	Į A

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

[#] cluttered chromatogram; sample peak coelutes with surrogate peak.

⁺The following descriptions of the TPH chromatogram are cursory in nature and McCampbell Analytical is not responsible for their interpretation: a) unmodified or weakly modified gasoline is significant; b) heavier gasoline range compounds are significant(aged gasoline?); c) lighter gasoline range compounds (the most mobile fraction) are significant; d) gasoline range compounds having broad chromatographic peaks are significant; biologically altered gasoline?; e) TPH pattern that does not appear to be derived from gasoline (stoddard solvent / mineral spirit?); f) one to a few isolated non-target peaks present; g) strongly aged gasoline or diesel range compounds are significant; h) lighter than water immiscible sheen/product is present; i) liquid sample that contains greater than ~2 vol. % sediment; j) reporting limit raised due to high MTBE content; k) TPH pattern that does not appear to be derived from gasoline (aviation gas). m) no recognizable pattern.

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P & D Environmental	Client Project ID: #0055; Former Haber	Date Sampled: 11/25/03
4020 Panama Court	Oil	Date Received: 11/26/03
Oakland, CA 94611-4931	Client Contact: Paul King	Date Extracted: 12/01/03-12/03/03
Odaland, CA 94011-4931	Client P.O.:	Date Analyzed: 12/01/03-12/03/03

	Chent 1.0.	•		Date	Analyzed. 12/01/05-1	2/03/0	13
Volatile Extraction Method: SW5030B	es Organics + Ox		•	P&T and GC/MS (Basic T	= '	Order: 0	311385
	· · · · · · · · · · · · · · · · · · ·	7.02	11,00011170				
Lab ID	<u> </u>			0311385-001B			
Client ID				MW 1	_ .		
Matrix				Water			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acetone	ND	1.0	5.0	tert-Amyl methyl ether (TAME)) ND	1.0	0.5
Benzene	2.5	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	1.0	t-Butyl alcohol (TBA)	ND	1.0	5.0
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	2-Chloroethyl Vinyl Ether	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	0.5
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
Ethanol	ND	1.0	50	Ethylbenzene	ND	1.0	0.5
Ethyl tert-butyl ether (ETBE)	ND	1.0	0.5	Hexachlorobutadiene	ND	1.0	0.5
2-Hexanone	ND	1.0	0.5	Iodomethane (Methyl iodide)	ND	1.0	5.0
Isopropylbenzene	ND	1.0	0.5	4-Isopropyl toluene	ND ND	1.0	0.5
Methanol	ND	1.0	500	Methyl-t-butyl ether (MTBE)	32	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 ND	1.0	0.5
1,1,2,2-Tetrachloroethane	ND	1.0	0.5	Tetrachloroethene	ND	1.0	0.5
Toluene	ND ND	1.0	0.5	1,2,3-Trichlorobenzene	ND	1.0	0.5
1,2,4-Trichlorobenzene	ND ND	1.0	0.5	1,1,1-Trichloroethane	ND	1.0	0.5
1,1,2-Trichloroethane	ND ND	1.0	0.5	Trichloroethene	ND ND	1.0	0.5
Trichlorofluoromethane	ND	1.0	0.5	1,2,3-Trichloropropane	ND	1.0	0.5
1,2,4-Trimethylbenzene	ND ND	1.0	0.5	1,3,5-Trimethylbenzene	ND	1.0	0.5
Vinyl Acetate	ND	1.0	5.0	Vinyl Chloride	ND ND	1.0	0.5
Xylenes	ND ND	1.0	0.5	v myr Chloride	ND i	1.0	. U.3
Ajtones	, 14D i			accurates (9/)			
0/001			ogate Ke	coveries (%)			
%SS1:	109	,		%SS2:	98.0		

%SS3: 98.4

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

ND means not detected above the reporting limit; N/A means analyte not applicable to this analysis.

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

h) lighter than water immiscible sheen/product is present; i) liquid sample that contains greater than ~2 vol. % sediment; j) sample diluted due to high organic content.



110 2nd Avenue South, #D7, Pacheco, CA 94553-5560
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P & D Environmental	Client Project ID: #0055; Former Haber	Date Sampled: 11/25/03
4020 Panama Court	Oil	Date Received: 11/26/03
Oakland, CA 94611-4931	Client Contact: Paul King	Date Extracted: 12/01/03-12/03/03
Oakiand, CA 94011-4931	Client P.O.:	Date Analyzed: 12/01/03-12/03/03

Volatiles Organics + Oxygenates by P&T and GC/MS (Basic Target List)*

Extraction Method: SW5030B Analytical Method: SW8260B Work Order: 0311385

Lab ID	İ			0311385-002B			
Client ID				MW 2			
Matrix				Water			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acetone	ND<2500	500	5.0	tert-Amyl methyl ether (TAME)	ND<250	500	0.5
Benzene	6800	500	0.5	Bromobenzene	ND<250	500	0.5
Bromochloromethane	ND<250	500	0.5	Bromodichloromethane	ND<250	500	0.5
Втоmoform	ND<250	500	0.5	Bromomethane	ND<250	500	0.5
2-Butanone (MEK)	ND<500	500	1.0	t-Butyl alcohol (TBA)	ND<2500	500	5.0
n-Butyl benzene	ND<250	500	0.5	sec-Butyl benzene	ND<250	500	0.5
tert-Butyl benzene	ND<250	500	0.5	Carbon Disulfide	ND<250	500	0.5
Carbon Tetrachloride	ND<250	500	0.5	Chlorobenzene	ND<250	500	0.5
Chloroethane	ND<250	500	0.5	2-Chloroethyl Vinyl Ether	ND<250	500	0.5
Chloroform	ND<250	500	0.5	Chloromethane	ND<250	500	0.5
2-Chlorotoluene	ND<250	500	0.5	4-Chlorotoluene	ND<250	500	0.5
Dibromochloromethane	ND<250	500	0.5	1,2-Dibromo-3-chloropropane	ND<250	500	0.5
1.2-Dibromoethane (EDB)	ND<250	500	0.5	Dibromomethane	ND<250	500	0.5
1,2-Dichlorobenzene	ND<250	500	0.5	1,3-Dichlorobenzene	ND<250	500	0.5
1.4-Dichlorobenzene	ND<250	500	0.5	Dichlorodifluoromethane	ND<250	500	0.5
1,1-Dichloroethane	ND<250	500	0.5	1,2-Dichloroethane (1,2-DCA)	ND<250	500	0.5
1.1-Dichloroethene	ND<250	500	0.5	cis-1,2-Dichloroethene	ND<250	500	0.5
trans-1,2-Dichloroethene	ND<250	500	0.5	1,2-Dichloropropane	ND<250	500	0.5
1,3-Dichloropropane	ND<250	500	0.5	2,2-Dichloropropane	ND<250	500	0.5
1,1-Dichloropropene	ND<250	500	0.5	cis-1,3-Dichloropropene	ND<250	500	0.5
trans-1,3-Dichloropropene	ND<250	500	0.5	Diisopropyl ether (DIPE)	ND<250	500	0.5
Ethanol	ND<25,000	500	50	Ethylbenzene	2900	500	0.5
Ethyl tert-butyl ether (ETBE)	ND<250	500	0.5	Hexachlorobutadiene	ND<250	500	0.5
2-Hexanone	ND<250	500	0.5	lodomethane (Methyl iodide)	ND<2500	500	5.0
Isopropylbenzene	ND<250	500	0.5	4-Isopropyl toluene	ND<250	500	0.5
Methanol	ND<250,000	500	500	Methyl-t-butyl ether (MTBE)	2700	500	0.5
Methylene chloride	ND<250	500	0.5	4-Methyl-2-pentanone (MIBK)	ND<250	500	0.5
Naphthalene	540	500	0.5	n-Propyl benzene	ND<250	500	0.5
Styrene	ND<250	500	0.5	1,1,1,2-Tetrachloroethane	ND<250	500	0.5
1,1,2,2-Tetrachloroethane	ND<250	500	0.5	Tetrachloroethene	ND<250	500	0.5
Toluene	8800	500	0.5	1,2,3-Trichlorobenzene	ND<250	500	0.5
1,2,4-Trichlorobenzene	ND<250	500	0.5	1.1.1-Trichloroethane	ND<250	500	0.5
1,1,2-Trichloroethane	ND<250	500	0.5	Trichloroethene	ND<250	500	0.5
Trichlorofluoromethane	ND<250	500	0.5	1,2,3-Trichloropropane	ND<250	500	0.5
1,2,4-Trimethylbenzene	1800	500	0.5	1,3,5-Trimethylbenzene	420	500	0.5
Vinyl Acetate	ND<2500	500	5.0	Vinyl Chloride	ND<250	500	0.5
Xylenes	16,000	500	0.5	. my · emoriae	1, 1, 2, 2, 0	200	. 0.5
1 Augustus 1	10,000	• • • • • • • •		ecoveries (%)			
%SS1:	10		IUgan N	%SS2:	99.3	<u> </u>	
	94.			/0334.	: 99.3	, 	· · · · · · · · · · · · · · · · · · ·
%SS3:	94.	1					

Comments:

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

ND means not detected above the reporting limit; N/A means analyte not applicable to this analysis.

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

h) lighter than water immiscible sheen/product is present; i) liquid sample that contains greater than ~2 vol. % sediment; j) sample diluted due to high organic content.

Jh

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P & D Environmental	Client Project ID: #0055; Former Haber	Date Sampled: 11/25/03
4020 Panama Court	Oil	Date Received: 11/26/03
Oakland, CA 94611-4931	Client Contact: Paul King	Date Extracted: 12/01/03-12/03/03
Oakiand, CA 94011-4931	Client P.O.:	Date Analyzed: 12/01/03-12/03/03

Volatiles Organics + Oxygenates by P&T and GC/MS (Basic Target List)*

Extraction Method: SW5030B Analytical Method: SW8260B Work Order: 0311385

Extraction Method: SW5030B		An	alytical Me	thod: SW8260B	Work	Order: 0	311385	
Lab ID		0311385-003B						
Client ID				MW 3				
Matrix				Water				
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit	
Acetone	ND<50	10	5.0	tert-Amyl methyl ether (TAME)	ND<5.0	10	0.5	
Benzene	ND<5.0	10	0.5	Bromobenzene	ND<5.0	10	0.5	
Bromochloromethane	ND<5.0	10	0.5	Bromodichloromethane	ND<5.0	10	0.5	
Bromoform	ND<5.0	10	0.5	Bromomethane	ND<5.0	10	0.5	
2-Butanone (MEK)	ND<10	10	1.0	t-Butyl alcohol (TBA)	ND<50	10	5.0	
n-Butyl benzene	ND<5.0	10	0.5	sec-Butyl benzene	ND<5.0	10	0.5	
tert-Butyl benzene	ND<5.0	10	0.5	Carbon Disulfide	ND<5.0	10	0.5	
Carbon Tetrachloride	ND<5.0	10	0.5	Chlorobenzene	ND<5.0	10	0.5	
Chloroethane	ND<5.0	10	0.5	2-Chloroethyl Vinyl Ether	ND<5.0	10	0.5	
Chloroform	ND<5.0	10	0.5	Chloromethane	ND<5.0	10	0.5	
2-Chlorotoluene	ND<5.0	10	0.5	4-Chlorotoluene	ND<5.0	10	0.5	
Dibromochloromethane	ND<5.0	10	0.5	1,2-Dibromo-3-chloropropane	ND<5.0	10	0.5	
1,2-Dibromoethane (EDB)	ND<5.0	10	0.5	Dibromomethane	: ND<5.0	10	0.5	
1,2-Dichlorobenzene	ND<5.0	10	0.5	1,3-Dichlorobenzene	ND<5.0	10	0.5	
1,4-Dichlorobenzene	ND<5.0	10	0.5	Dichlorodifluoromethane	ND<5.0	10	0.5	
1,1-Dichloroethane	ND<5.0	10	0.5	1,2-Dichloroethane (1,2-DCA)	ND<5.0	10	0.5	
1,1-Dichloroethene	ND<5.0	10	0.5	cis-1,2-Dichloroethene	ND<5.0	10	0.5	
trans-1,2-Dichloroethene	ND<5.0	10	0.5	1,2-Dichloropropane	ND<5.0	10	0.5	
1,3-Dichloropropane	ND<5.0	10	0.5	2,2-Dichloropropane	ND<5.0	10	0.5	
1,1-Dichloropropene	ND<5.0	10	0.5	cis-1,3-Dichloropropene	ND<5.0	10	0.5	
trans-1,3-Dichloropropene	ND<5.0	10	0.5	Diisopropyl ether (DIPE)	ND<5.0	10	0.5	
Ethanol	ND<500	10	50	Ethylbenzene	ND<5.0	10	0.5	
Ethyl tert-butyl ether (ETBE)	ND<5.0	10	0.5	Hexachlorobutadiene	ND<5.0	10	0.5	
2-Hexanone	ND<5.0	10	0.5	Iodomethane (Methyl iodide)	ND<50	10	5.0	
Isopropylbenzene	ND<5.0	10	0.5	4-Isopropyl toluene	ND<5.0	10	0.5	
Methanol	ND<5000	10	500	Methyl-t-butyl ether (MTBE)	330	10	0.5	
Methylene chloride	ND<5.0	10	0.5	4-Methyl-2-pentanone (MIBK)	ND<5.0	10	0.5	
Naphthalene	ND<5.0	10	0.5	n-Propyl benzene	ND<5.0	10	0.5	
Styrene	ND<5.0	10	0.5	1,1,1,2-Tetrachloroethane	ND<5.0	10	0.5	
1,1,2,2-Tetrachloroethane	ND<5.0	10	0.5	Tetrachloroethene	ND<5.0	10	0.5	
Toluene	ND<5.0	10	0.5	1.2.3-Trichlorobenzene	ND<5.0	10	0.5	
1,2,4-Trichlorobenzene	ND<5.0	10	0.5	1,1,1-Trichloroethane	ND<5.0	10	0.5	
1.1.2-Trichloroethane	ND<5.0	10	0.5	Trichloroethene	ND<5.0	10	0.5	
Trichlorofluoromethane	ND<5.0	10	0.5	1,2,3-Trichloropropane	ND<5.0	10	0.5	
1,2,4-Trimethylbenzene	ND<5.0	10	0.5	1,3,5-Trimethylbenzene	ND<5.0	10	0.5	
Vinyl Acetate	ND<50	10	5.0	Vinyl Chloride	ND<5.0	10	0.5	
Xylenes	ND<5.0	10	0.5	i myt Omorau	1427-7-10		1. 0.5	
21 tolled	1422-240			ecoveries (%)				
%SS1:	106		. ogate M	%SS2:	101			
%\$\$3:	94.4				101			
Comments	7							

Comments

Angela Rydelius, Lab Manager

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

ND means not detected above the reporting limit; N/A means analyte not applicable to this analysis.

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

h) lighter than water immiscible sheen/product is present; i) liquid sample that contains greater than ~2 vol. % sediment; j) sample diluted due to high organic content.

110 2nd Avenue South, #D7, Pacheco, CA 94553-5560
Telephone: 925-798-1620 Fax: 925-798-1622
http://www.mccampbell.com E-mail: main@mccampbell.com

P & D Environmental	Client Project ID: #0055; Former Haber Oil Client Contact: Paul King Client P.O.:	Date Sampled: 11/25/03
4020 Panama Court	Oti	Date Received: 11/26/03
Oakland, CA 94611-4931	Client Contact: Paul King	Date Extracted: 12/01/03-12/03/03
Oakiaiu, CA 94011-4931	Client P.O.:	Date Analyzed: 12/01/03-12/03/03

Volatiles Organics + Oxygenates by P&T and GC/MS (Basic Target List)*

Extraction Method: SW 5030B Analytical Method: SW 8260B Work Order: 0311385

 Lab ID
 0311385-004B

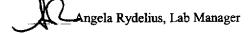
 Client ID
 MW 4

 Matrix
 Water

Matrix Water							
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acetone	ND<2500	500	5.0	tert-Amyl methyl ether (TAME)	ND<250	500	0.5
Benzene	ND<250	500	0.5	Вготовепие	ND<250	500	0.5
Bromochloromethane	ND<250	500	0.5	Bromodichloromethane	ND<250	500	0.5
Bromoform	ND<250	500	0.5	Bromomethane	ND<250	500	0.5
2-Butanone (MEK)	ND<500	500	1.0	t-Butyl alcohol (TBA)	ND<2500	500	5.0
n-Butyl benzene	ND<250	500	0.5	sec-Butyl benzene	ND<250	500	0.5
tert-Butyl benzene	ND<250	500	0.5	Carbon Disulfide	ND<250	500	0.5
Carbon Tetrachloride	ND<250	500	0.5	Chlorobenzene	ND<250	500	0.5
Chloroethane	ND<250	500	0.5	2-Chloroethyl Vinyl Ether	ND<250	500	0.5
Chloroform	ND<250	500	0.5	Chloromethane	ND<250	500	0.5
2-Chlorotoluene	ND<250	500	0.5	4-Chlorotoluene	ND<250	500	0.5
Dibromochloromethane	ND<250	500	0.5	1,2-Dibromo-3-chloropropane	ND<250	500	0.5
1,2-Dibromoethane (EDB)	ND<250	500	0.5	Dibromomethane	ND<250	500	0.5
1,2-Dichlorobenzene	ND<250	500	0.5	1,3-Dichlorobenzene	ND<250	500	0.5
1,4-Dichlorobenzene	ND<250	500	0.5	Dichlorodifluoromethane	ND<250	500	0.5
1,1-Dichloroethane	ND<250	500	0.5	1,2-Dichloroethane (1,2-DCA)	ND<250	500	0.5
1,1-Dichloroethene	ND<250	500	0.5	cis-1,2-Dichloroethene	ND<250	500	0.5
trans-1,2-Dichloroethene	ND<250	500	0.5	1,2-Dichloropropane	ND<250	500	0.5
1,3-Dichloropropane	ND<250	500	0.5	2,2-Dichloropropane	ND<250	500	0.5
1,1-Dichloropropene	ND<250	500		cis-1,3-Dichloropropene	ND<250	500	0.5
trans-1,3-Dichloropropene	ND<250	500	0.5	Diisopropyl ether (DIPE)	ND<250	500	0.5
Ethanol	ND<25,000	500	50	Ethylbenzene	ND<250	500	0.5
Ethyl tert-butyl ether (ETBE)	ND<250	500	0.5	Hexachlorobutadiene	ND<250	500	0.5
2-Hexanone	ND<250	500	0.5	Iodomethane (Methyl iodide)	ND<2500	500	5.0
Isopropylbenzene	ND<250	500	0.5	4-Isopropyl toluene	ND<250	500	0.5
Methanol	ND<250,000	500	500	Methyl-t-butyl ether (MTBE)	8800	500	0.5
Methylene chloride	ND<250	500	0.5	4-Methyl-2-pentanone (MIBK)	ND<250	500	0.5
Naphthalene	ND<250	500	0.5	n-Propyl benzene	ND<250	500	0.5
Styrene	ND<250	500	0.5	1,1,1,2-Tetrachloroethane	ND<250	500	0.5
1,1,2,2-Tetrachloroethane	ND<250	500	0.5	Tetrachloroethene	ND<250	500	0.5
Toluene	ND<250	500	0.5	1,2,3-Trichlorobenzene	ND<250	500	0.5
1,2,4-Trichlorobenzene	ND<250	500	0.5	1,1,1-Trichloroethane	ND<250	500	0.5
1,1,2-Trichloroethane	ND<250	500	0.5	Trichloroethene	ND<250	500	0.5
Trichlorofluoromethane	ND<250	500	0.5	1,2,3-Trichloropropane	ND<250	500	0.5
1,2,4-Trimethylbenzene	ND<250	500	0.5	1,3,5-Trimethylbenzene	ND<250	500	0.5
Vinyl Acetate	ND<2500	500	5.0	Vinyl Chloride	ND<250	500	0.5
Xylenes	ND<250	500	0.5				
		Sur	rogate R	ecoveries (%)			
%SS1:	10			%SS2:	101		
%SS3:	92.	9					

Comments:

h) lighter than water immiscible sheen/product is present; i) liquid sample that contains greater than ~2 vol. % sediment; j) sample diluted due to high organic content.



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

ND means not detected above the reporting limit; N/A means analyte not applicable to this analysis.

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

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P & D Environmental	Client Project ID: #0055; Former Haber	Date Sampled: 11/25/03
4020 Panama Court	Oil	Date Received: 11/26/03
Oakland, CA 94611-4931	Client Contact: Paul King	Date Extracted: 12/01/03-12/03/03
Oakiand, CA 94011-4931	Client P.O.:	Date Analyzed: 12/01/03-12/03/03

Volatiles Organics + Oxygenates by P&T and GC/MS (Basic Target List)*

Extraction Method: SW5030B	Analytical Method: SW8260B	Work Order: 0311385

Lab ID				0311385-005B					
Client ID				MW 5					
Matrix	Matrix Water								
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reportin Limit		
Acetone	ND	1.0	5.0	tert-Amyl methyl ether (TAME)	ND	1.0	0.5		
Вепzепе	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	1.0	t-Butyl alcohol (TBA)	ND	1.0	5.0		
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	2-Chloroethyl Vinyl Ether	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	0.5		
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		
Ethanol	ND	1.0	50	Ethylbenzene	ND	1.0	0.5		
Ethyl tert-butyl ether (ETBE)	ND	1.0	0.5	Hexachlorobutadiene	ND	1.0	0.5		
2-Hexanone	ND	1.0	0.5	lodomethane (Methyl iodide)	ND	1.0	5.0		
Isopropylbenzene	ND	1.0	0.5	4-Isopropyl toluene	ND	1.0	0.5		
Methanol	ND	1.0	500	Methyl-t-butyl ether (MTBE)	0.84	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		1,2,3-Trichlorobenzene	ND	1.0	0.5		
1,2,4-Trichlorobenzene	ND	1.0		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	ND	1.0	0.5	1,3,5-Trimethylbenzene	ND	1.0	0.5		
Vinyl Acetate	ND	1.0	5.0	Vinyl Chloride	ND	1.0	0.5		
Xylenes	ND	1.0	0.5	-	1		· · · · ·		
		Sur	_	ecoveries (%)	~~~	•			
%SS1:	106	_		%SS2:	101				
%SS3:	95.8		, , , , , , , , , , , , , , , , , , , ,	101					

Jh.

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

ND means not detected above the reporting limit; N/A means analyte not applicable to this analysis.

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

h) lighter than water immiscible sheen/product is present; i) liquid sample that contains greater than -2 vol. % sediment; j) sample diluted due to high organic content.

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P & D Environmental	Client Project ID: #0055; Former Haber	Date Sampled: 11/25/03
4020 Panama Court	Oil	Date Received: 11/26/03
Oakland, CA 94611-4931	Client Contact: Paul King	Date Extracted: 12/01/03-12/03/03
Oliciand, Cri 94011-4551	Client P.O.;	Date Analyzed: 12/01/03-12/03/03

Volatiles Organics + Oxygenates by P&T and GC/MS (Basic Target List)*

 Extraction Method:
 SW5030B
 Analytical Method:
 SW8260B
 Work Order:
 0311385

 Lab ID
 0311385-006B
 O3100 MW 6
Compound Concentration DF Limb Compound Concentration DF	Client ID	MW 6							
Compound Concentration DF Lient Compound Concentration DF	Matrix				Water				
Benzene ND 1.0 0.5 Bromobenzene ND 1.0 1.0 Bromochloromethane ND 1.0 0.5 Bromodichloromethane ND 1.0 1.0 1.0 Bromochloromethane ND 1.0 1.0 1.0 Bromoform ND 1.0 I.0 I	Compound	Concentration *	DF		Compound	Concentration *	DF	Reporting Limit	
Bromochloromethane	Acetone	ND	1.0	5.0	tert-Amyl methyl ether (TAME)	ND	1.0	0.5	
Bromoform	Benzene	ND	1.0	0.5	Bromobenzene	ND	1.0	0.5	
2-Butanone (MEK) ND 1.0 1-Butyl alcohol (TBA) ND 1.0 N-Butyl benzene ND 1.0 0.5 sec-Butyl benzene ND 1.0 0.5 Carbon Disulfide ND 1.0 0.5 Chlorocethane ND 1.0 0.5 Chloroce	Bromochloromethane	ND	1.0	0.5	Bromodichloromethane	ND	1.0	0.5	
n-Butyl benzene		ND	1.0	0.5	Bromomethane	ND	1.0	0.5	
tert-Butyl benzene ND 1.0 0.5 Carbon Disulfide ND 1.0 Carbon Tetrachloride ND 1.0 0.5 Chlorobenzene ND 1.0 Chloroethane ND 1.0 0.5 2-Chloroethyl Vinyl Ether ND 1.0 Chloroform 0.89 1.0 0.5 Chloromethane ND 1.0 2-Chlorotoluene ND 1.0 0.5 4-Chlorotoluene ND 1.0 Dibromochloromethane ND 1.0 0.5 1,2-Dibromo-3-chloropropane ND 1.0 1,2-Dibromochloromethane ND 1.0 0.5 1,2-Dibromo-3-chloropropane ND 1.0 1,2-Dichlorosthane ND 1.0 0.5 1,3-Dichloropropane ND 1.0 1,1-Dichlorosthane ND 1.0 0.5 1,2-Dichlorosthene ND 1.0 1,1-Dichloroptopene ND 1.0 0.5 1,2-Dichloroptopene ND 1.0 trans-1,2-Dichloroptopene ND 1.0	2-Butanone (MEK)	ND	1.0	1.0	t-Butyl alcohol (TBA)	ND	1.0	5.0	
Carbon Tetrachloride ND 1.0 0.5 Chloroethare ND 1.0 Chloroethane ND 1.0 0.5 2-Chloroethyl Vinyl Ether ND 1.0 Chloroform 0.89 1.0 0.5 2-Chlorotoluene ND 1.0 2-Chlorotoluene ND 1.0 0.5 4-Chlorotoluene ND 1.0 1,2-Dibromochloromethane ND 1.0 0.5 1,2-Dibromo-3-chloropropane ND 1.0 1,2-Dibromochloromethane ND 1.0 0.5 Dibromomethane ND 1.0 1,2-Dibromochane (EDB) ND 1.0 0.5 Dibromomethane ND 1.0 1,2-Dichloroethane (EDB) ND 1.0 0.5 Dichlorodifluoromethane ND 1.0 1,1-Dichloroethane ND 1.0 0.5 1,2-Dichloroethane ND 1.0 1,1-Dichloroethane ND 1.0 0.5 1,2-Dichloroethane ND 1.0 1,1-Dichloroethane ND 1.0	n-Butyl benzene	ND	1.0	0.5	sec-Butyl benzene	ND	1.0	: 0.5	
Chloroethane ND 1.0 0.5 2-Chloroethyl Vinyl Ether ND 1.0 Chloroform 0.89 1.0 0.5 Chloromethane ND 1.0 2-Chlorotoluene ND 1.0 0.5 Chlorotoluene ND 1.0 Dibromochloromethane ND 1.0 0.5 1,2-Dibromo-3-chloropropane ND 1.0 1,2-Dibromochloromethane ND 1.0 0.5 1,2-Dibromo-3-chloropropane ND 1.0 1,2-Dichlorobenzene ND 1.0 0.5 1,3-Dichlorobenzene ND 1.0 1,4-Dichlorobenzene ND 1.0 0.5 1,2-Dichlorodifluoromethane ND 1.0 1,4-Dichlorobenzene ND 1.0 0.5 1,2-Dichlorofifluoromethane ND 1.0 1,1-Dichlorobethane ND 1.0 0.5 1,2-Dichloroethane ND 1.0 1,1-Dichlorobethane ND 1.0 0.5 1,2-Dichloroethane ND 1.0 1,1-Dichlorobethane ND	tert-Butyl benzene	ND	1.0	0.5	Carbon Disulfide	ND	1.0	0.5	
Chloroform	Carbon Tetrachloride	ND	1.0	0.5	Chlorobenzene	ND	1.0	0.5	
2-Chlorotoluene	Chloroethane	ND	1.0	0.5	2-Chloroethyl Vinyl Ether	ND	1.0	0.5	
Dibromochloromethane ND	Chloroform	0.89	1.0	0.5	Chloromethane	ND	1.0	0.5	
1,2-Dibromoethane (EDB) ND 1.0 0.5 Dibromomethane ND 1.0 1,2-Dichlorobenzene ND 1.0 0.5 1,3-Dichlorobenzene ND 1.0 1,4-Dichlorobenzene ND 1.0 0.5 Dichlorodifluoromethane ND 1.0 1,1-Dichloroethane ND 1.0 0.5 1,2-Dichloroethane (1,2-DCA) ND 1.0 1,1-Dichloroethene ND 1.0 0.5 1,2-Dichloroethene ND 1.0 1,1-Dichloroethene ND 1.0 0.5 1,2-Dichloropropane ND 1.0 1,3-Dichloropropane ND 1.0 0.5 2,2-Dichloropropane ND 1.0 1,1-Dichloropropane ND 1.0 0.5 2,2-Dichloropropane ND 1.0 1,1-Dichloropropane ND 1.0 0.5 2,2-Dichloropropane ND 1.0 1,1-Dichloropropane ND 1.0 0.5 bisopropyle ther (DIPE) ND 1.0 1,1-Dichloropropane ND	2-Chlorotoluene	ND	1.0	0.5	4-Chlorotoluene	ND	1.0	0.5	
1,2-Dibromoethane (EDB) ND 1.0 0.5 Dibromomethane ND 1.0 1,2-Dichlorobenzene ND 1.0 0.5 1,3-Dichlorobenzene ND 1.0	Dibromochloromethane	ND	1.0	0.5	1,2-Dibromo-3-chloropropane	ND	1.0	0.5	
1,4-Dichlorobenzene ND 1.0 0.5 Dichlorodifluoromethane ND 1.0 1,1-Dichloroethane ND 1.0 0.5 1,2-Dichloroethane (1,2-DCA) ND 1.0 1,1-Dichloroethene ND 1.0 0.5 1,2-Dichloroethene ND 1.0 1,3-Dichloroethene ND 1.0 0.5 1,2-Dichloropropane ND 1.0 1,3-Dichloropropane ND 1.0 0.5 cis-1,3-Dichloropropane ND 1.0 1,1-Dichloropropene ND 1.0 0.5 cis-1,3-Dichloropropene ND 1.0 1,1-Dichloropropene ND 1.0 0.5 Diisopropyl ether (DIPE) ND 1.0 1,1-Dichloropropene ND 1.0 0.5 Ethylenere ND 1.0 1,1-Dichloropropene ND 1.0 0.5 Diisopropyle ther (DIPE) ND 1.0 1,1-Dichloropropene ND 1.0 0.5 Ethylenere ND 1.0 1,0 1.0 1.0	1,2-Dibromoethane (EDB)	ND	1.0	0.5				0.5	
1,4-Dichlorobenzene ND 1.0 0.5 Dichlorodifluoromethane ND 1.0 1,1-Dichloroethane ND 1.0 0.5 1,2-Dichloroethane (1,2-DCA) ND 1.0 1,1-Dichloroethene ND 1.0 0.5 1,2-Dichloroethene ND 1.0 1,3-Dichloroethene ND 1.0 0.5 1,2-Dichloropropane ND 1.0 1,3-Dichloropropane ND 1.0 0.5 cis-1,3-Dichloropropane ND 1.0 1,1-Dichloropropene ND 1.0 0.5 cis-1,3-Dichloropropene ND 1.0 1,1-Dichloropropene ND 1.0 0.5 Diisopropyl ether (DIPE) ND 1.0 1,1-Dichloropropene ND 1.0 0.5 Ethylenere ND 1.0 1,1-Dichloropropene ND 1.0 0.5 Diisopropyle ther (DIPE) ND 1.0 1,1-Dichloropropene ND 1.0 0.5 Ethylenere ND 1.0 1,0 1.0 1.0	1,2-Dichlorobenzene	ND	1.0	0.5	1,3-Dichlorobenzene	ND	1.0	0.5	
1,1-Dichloroethene	1,4-Dichlorobenzene	ND	1.0	0.5		ND	1.0	0.5	
1,1-Dichloroethene ND 1.0 0.5 cis-1,2-Dichloroethene ND 1.0 1,3-Dichloroethene ND 1.0 0.5 1,2-Dichloropropane ND 1.0 1,3-Dichloropropane ND 1.0 0.5 2,2-Dichloropropane ND 1.0 1,1-Dichloropropene ND 1.0 0.5 cis-1,3-Dichloropropene ND 1.0 trans-1,3-Dichloropropene ND 1.0 0.5 Diisopropyl ether (DIPE) ND 1.0 Ethanol ND 1.0 50 Ethylbenzene ND 1.0 Ethyl tert-butyl ether (ETBE) ND 1.0 0.5 Hexachlorobutadiene ND 1.0 2-Hexanone ND 1.0 0.5 Hexachlorobutadiene ND 1.0 1-Sopropylbenzene ND 1.0 0.5 Hexachlorobutadiene ND 1.0 1-Sopropylbenzene ND 1.0 0.5 Hexachlorobutaene ND 1.0 Methyler-butylenzene ND 1.0	1,1-Dichloroethane	ND	1.0	0.5	1,2-Dichloroethane (1,2-DCA)	ND	1.0	0.5	
trans-1,2-Dichloroethene ND 1.0 0.5 1,2-Dichloropropane ND 1.0 1,3-Dichloropropane ND 1.0 0.5 2,2-Dichloropropane ND 1.0 1,1-Dichloropropene ND 1.0 0.5 cis-1,3-Dichloropropene ND 1.0 trans-1,3-Dichloropropene ND 1.0 0.5 Diisopropyl ether (DIPE) ND 1.0 Ethal collection ND 1.0 0.5 Ethylbenzene ND 1.0 Ethyl tert-butyl ether (ETBE) ND 1.0 0.5 Hexachlorobutadiene ND 1.0 2-Hexanone ND 1.0 0.5 Hexachlorobutadiene ND 1.0 Isopropylbenzene ND 1.0 0.5 Hexachlorobutadiene ND 1.0 Methanol ND 1.0 0.5 Jelsopropyl toluene ND 1.0 Methanol ND 1.0 0.5 Jelsopropyl toluene ND 1.0 Methanol ND 1.0 0.5	1,1-Dichloroethene	ND	1.0	0.5	cis-1.2-Dichloroethene		1.0	0.5	
1,3-Dichloropropane ND 1.0 0.5 2,2-Dichloropropane ND 1.0 1,1-Dichloropropene ND 1.0 0.5 cis-1,3-Dichloropropene ND 1.0 trans-1,3-Dichloropropene ND 1.0 0.5 Diisopropyl ether (DIPE) ND 1.0 Ethanol ND 1.0 50 Ethylbenzene ND 1.0 Ethyl tert-butyl ether (ETBE) ND 1.0 0.5 Hexachlorobutadiene ND 1.0 2-Hexanone ND 1.0 0.5 Hexachlorobutadiene ND 1.0 Isopropylbenzene ND 1.0 0.5 Hexachlorobutadiene ND 1.0 Isopropylbenzene ND 1.0 0.5 Hexachlorobutadiene ND 1.0 Methylene chloride ND 1.0 0.5 Hexachloropyl toluene ND 1.0 Methylene chloride ND 1.0 0.5 4-Methyl-2-pentanone (MIBK) ND 1.0 Naphthalene ND 1.0	trans-1,2-Dichloroethene		1.0	0.5		· · · · · · · · · · · · · · · · · · ·		0.5	
1,1-Dichloropropene ND 1.0 0.5 cis-1,3-Dichloropropene ND 1.0 trans-1,3-Dichloropropene ND 1.0 0.5 Diisopropyl ether (DIPE) ND 1.0 Ethanol ND 1.0 50 Ethylbenzene ND 1.0 Ethyl tert-butyl ether (ETBE) ND 1.0 0.5 Hexachlorobutadiene ND 1.0 2-Hexanone ND 1.0 0.5 Hexachlorobutadiene ND 1.0 Isopropylbenzene ND 1.0 0.5 Hexachlorobutadiene ND 1.0 Methylene chloride ND 1.0 0.5 4-Isopropyl toluene ND 1.0 Methylene chloride ND 1.0 0.5 4-Methyl-2-pentanone (MIBK) ND 1.0 Methylene chloride ND 1.0 0.5 1-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 1-I,1,2-Tetrachloroethane ND 1.0 1,1,2-Tetrachloroethane ND 1.0	1,3-Dichloropropane			1				0.5	
trans-1,3-Dichloropropene ND 1.0 0.5 Diisopropyl ether (DIPE) ND 1.0 Ethanol ND 1.0 50 Ethylbenzene ND 1.0 Ethyl tert-butyl ether (ETBE) ND 1.0 0.5 Hexachlorobutadiene ND 1.0 2-Hexanone ND 1.0 0.5 Idodmethane (Methyl iodide) ND 1.0 Isopropylbenzene ND 1.0 0.5 Idomethane (Methyl iodide) ND 1.0 Methanol ND 1.0 0.5 4-Isopropyl toluene ND 1.0 Methylene chloride ND 1.0 500 Methyl-t-butyl ether (MTBE) ND 1.0 Methylene chloride ND 1.0 0.5 4-Methyl-2-pentanone (MIBK) ND 1.0 Methylene chloride ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5	1,1-Dichloropropene	ND	1.0	0.5		· · · · · · · · · · · · · · · · · · ·		0.5	
Ethanol ND 1.0 50 Ethylbenzene ND 1.0 Ethyl tert-butyl ether (ETBE) ND 1.0 0.5 Hexachlorobutadiene ND 1.0 2-Hexanone ND 1.0 0.5 Iodomethane (Methyl iodide) ND 1.0 Isopropylbenzene ND 1.0 0.5 4-Isopropyl toluene ND 1.0 Methanol ND 1.0 500 Methyl-t-butyl ether (MTBE) ND 1.0 Methylene chloride ND 1.0 0.5 4-Methyl-2-pentanone (MIBK) ND 1.0 Naphthalene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 1,1,1,2-Tetrachloroethane ND 1.0 1,1,2,2-Tetrachloroethane ND 1.0 0.5 Tetrachloroethene ND 1.0 1,2,4-Trichloroethane ND 1.0 0.5	trans-1,3-Dichloropropene	ND	1.0	0.5				0.5	
Ethyl tert-butyl ether (ETBE) ND 1.0 0.5 Hexachlorobutadiene ND 1.0 2-Hexanone ND 1.0 0.5 Iodomethane (Methyl iodide) ND 1.0 Isopropylbenzene ND 1.0 0.5 4-Isopropyl toluene ND 1.0 Methylene chloride ND 1.0 500 Methyl-t-butyl ether (MTBE) ND 1.0 Methylene chloride ND 1.0 0.5 4-Methyl-2-pentanone (MIBK) ND 1.0 Naphthalene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 1,1,1,2-Tetrachloroethane ND 1.0 1,1,2,2-Tetrachloroethane ND 1.0 0.5 Tetrachloroethene ND 1.0 1,2,4-Trichloroethane ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 1,2,4-Trimethylbenzene ND 1.0 <td></td> <td>ND</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>0.5</td>		ND						0.5	
2-Hexanone ND 1.0 0.5 Iodomethane (Methyl iodide) ND 1.0 Isopropylbenzene ND 1.0 0.5 4-Isopropyl toluene ND 1.0 Methanol ND 1.0 500 Methyl-t-butyl ether (MTBE) ND 1.0 Methylene chloride ND 1.0 0.5 4-Methyl-2-pentanone (MIBK) ND 1.0 Naphthalene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 1,1,1,2-Tetrachloroethane ND 1.0 1,1,2,2-Tetrachloroethane ND 1.0 0.5 Tetrachloroethene ND 1.0 1,2,4-Trichlorobenzene ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 <td>Ethyl tert-butyl ether (ETBE)</td> <td>+</td> <td>1.0</td> <td></td> <td>_ · · · · · · · · · · · · · · · · · · ·</td> <td></td> <td></td> <td>0.5</td>	Ethyl tert-butyl ether (ETBE)	+	1.0		_ · · · · · · · · · · · · · · · · · · ·			0.5	
Isopropylbenzene ND 1.0 0.5 4-Isopropyl toluene ND 1.0 Methanol ND 1.0 500 Methyl-t-butyl ether (MTBE) ND 1.0 Methylene chloride ND 1.0 0.5 4-Methyl-2-pentanone (MIBK) ND 1.0 Naphthalene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 1,1,1,2-Tetrachloroethane ND 1.0 1,1,2,2-Tetrachloroethane ND 1.0 0.5 Tetrachloroethene ND 1.0 Toluene ND 1.0 0.5 1,2,3-Trichloroethene ND 1.0 1,2,4-Trichloroethane ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 Trichlorofluoromethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 Trichlorofluoromethane ND 1.0		ND	1.0	0.5		<u> </u>		5.0	
Methanol ND 1.0 500 Methyl-t-butyl ether (MTBE) ND 1.0 Methylene chloride ND 1.0 0.5 4-Methyl-2-pentanone (MIBK) ND 1.0 Naphthalene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 1,1,1,2-Tetrachloroethane ND 1.0 1,1,2,2-Tetrachloroethane ND 1.0 0.5 Tetrachloroethene ND 1.0 Toluene ND 1.0 0.5 1,2,3-Trichlorobenzene ND 1.0 1,2,4-Trichlorobenzene ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 Tylenethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 <td< td=""><td>Isopropylbenzene</td><td>ND</td><td>1.0</td><td></td><td></td><td>· · · · · · · · · · · · · · · · · · ·</td><td></td><td>0.5</td></td<>	Isopropylbenzene	ND	1.0			· · · · · · · · · · · · · · · · · · ·		0.5	
Methylene chloride ND 1.0 0.5 4-Methyl-2-pentanone (MIBK) ND 1.0 Naphthalene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 1,1,1,2-Tetrachloroethane ND 1.0 1,1,2,2-Tetrachloroethane ND 1.0 0.5 Tetrachloroethane ND 1.0 Toluene ND 1.0 0.5 1,2,3-Trichlorobenzene ND 1.0 1,2,4-Trichlorobenzene ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 1,2,4-Trimethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0	Methanol	ND				·		0.5	
Naphthalene ND 1.0 0.5 n-Propyl benzene ND 1.0 Styrene ND 1.0 0.5 1,1,1,2-Tetrachloroethane ND 1.0 1,1,2,2-Tetrachloroethane ND 1.0 0.5 Tetrachloroethane ND 1.0 Toluene ND 1.0 0.5 1,2,3-Trichlorobenzene ND 1.0 1,2,4-Trichloroethane ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 1,2,4-Trimethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0	Methylene chloride	ND				- 		0.5	
Styrene ND 1.0 0.5 1,1,1,2-Tetrachloroethane ND 1.0 1,1,2,2-Tetrachloroethane ND 1.0 0.5 Tetrachloroethane ND 1.0 Toluene ND 1.0 0.5 1,2,3-Trichloroethane ND 1.0 1,2,4-Trichloroethane ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 1,2,4-Trimethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0	Naphthalene	ND		0.5				0.5	
1,1,2,2-Tetrachloroethane ND 1.0 0.5 Tetrachloroethene ND 1.0 Toluene ND 1.0 0.5 1,2,3-Trichlorobenzene ND 1.0 1,2,4-Trichlorobenzene ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 1,2,4-Trimethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0	Styrene			0.5		·+		0.5	
Toluene ND 1.0 0.5 1,2,3-Trichlorobenzene ND 1.0 1,2,4-Trichlorobenzene ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 1,2,4-Trimethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0	•		1.0	+				0.5	
1,2,4-Trichlorobenzene ND 1.0 0.5 1,1,1-Trichloroethane ND 1.0 1,1,2-Trichloroethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 1,2,4-Trimethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0	Toluene	ND						0.5	
1,1,2-Trichloroethane ND 1.0 0.5 Trichloroethene ND 1.0 Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 1,2,4-Trimethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0				• • • • • • • • • • • • • • • • • • • 				0.5	
Trichlorofluoromethane ND 1.0 0.5 1,2,3-Trichloropropane ND 1.0 1,2,4-Trimethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0						· · · · · · · · · · · · · · · · · · ·	<u>_</u> _	0.5	
1,2,4-Trimethylbenzene ND 1.0 0.5 1,3,5-Trimethylbenzene ND 1.0 Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0				 				0.5	
Vinyl Acetate ND 1.0 5.0 Vinyl Chloride ND 1.0								0.5	
						+		0.5	
	Xylenes	ND		0.5		-1	1.0	.1 0.5	
Surrogate Recoveries (%)					coveries (%)				
%SS1: 112 %SS2: 103	%SS1:	112				103			
%SS3: 103	%SS3:	103	1						

Comments

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

ND means not detected above the reporting limit; N/A means analyte not applicable to this analysis.

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

h) lighter than water immiscible sheen/product is present; i) liquid sample that contains greater than ~2 vol. % sediment; j) sample diluted due to high organic content.

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P & D Environmental	Client Project ID: #0055; Former Haber	Date Sampled: 11/25/03				
4020 Panama Court	Oil	Date Received: 11/26/03				
Oakland, CA 94611-4931	Client Contact: Paul King	Date Extracted: 12/01/03-12/02/03				
Oakland, CA 94011-4931	Client P.O.:	Date Analyzed: 12/01/03-12/02/03				

Volatiles Organics + Oxygenates by P&T and GC/MS (Basic Target List)*

 Extraction Method:
 SW5030B
 Analytical Method:
 SW8260B
 Work Order:
 0311385

 Lab ID
 0311385-007B

Labil				0311385-007B					
Client II)	MW 7							
Matrix	(Water							
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit		
Acetone	ND	1.0	5.0	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	1.0	t-Butyl alcohol (TBA)	ND	1.0	5.0		
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	2-Chloroethyl Vinyl Ether	ND	1.0	0.5		
Chloroform	0.76	1.0	0.5	Chloromethane	ND	1.0	0.5		
2-Chlorotoluene	ND	1.0	0.5	4-Chiorotoluene	ND	1.0	0.5		
Dibromochloromethane	ND	1.0	0.5	1,2-Dibromo-3-chloropropane	ND	1.0	0.5		
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		
Ethanol	ND	1.0	50	Ethylbenzene	ND	1.0	0.5		
Ethyl tert-butyl ether (ETBE)	ND	1.0	0.5	Hexachlorobutadiene	ND	1.0	0.5		
2-Hexanone	ND	1.0	0.5	Iodomethane (Methyl iodide)	ND	1.0	5.0		
Isopropylbenzene	ND	0.1	0.5	4-Isopropyl toluene	ND	1.0	0.5		
Methanol	ND	1.0	500	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	0.78	1.0	0.5		
Toluene	ND	1.0	0.5	1,2,3-Trichlorobenzene	ND 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	ND	1.0	0.5	1,3,5-Trimethylbenzene	ND	1.0	0.5		
Vinyl Acetate	ND	1.0	5.0	Vinyl Chloride	ND	1.0	0.5		
Xylenes	ND	1.0	0.5		<u></u>				
				ecoveries (%)					
%SS1:	111 %SS2: 99.7								
%SS3:	99.			,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	75.				
/0000.	99.	v		1					

Comments:

J

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

ND means not detected above the reporting limit; N/A means analyte not applicable to this analysis.

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

h) lighter than water immiscible sheen/product is present; i) liquid sample that contains greater than ~2 vol. % sediment; j) sample diluted due to high organic content.

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P & D Environmental	Client Project ID: #0055; Former Haber	Date Sampled: 11/25/03				
4020 Panama Court	Oil	Date Received: 11/26/03				
Oakland, CA 94611-4931	Client Contact: Paul King	Date Extracted: 12/01/03-12/03/03				
Oakimit, C/1 7-011-4931	Client P.O.:	Date Analyzed: 12/01/03-12/03/03				

Volatiles Organics + Oxygenates by P&T and GC/MS (Basic Target List)*

Extraction Method: SW5030B Analytical Method: SW8260B Work Order: 0311385

			ary dour 1110	alou. 5W 8200D	WOIR	Order, o	311302
Lab ID				0311385-008B			
Client ID				MW 8			
Matrix				Water			
Compound	Concentration *	DF	Reporting Limit	Compound	Concentration *	DF	Reporting Limit
Acetone	ND	1.0	5.0	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	1.0	t-Butyl alcohol (TBA)	ND	1.0	5.0
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	2-Chloroethyl Vinyl Ether	ND	1.0	0.5
Chloroform	1.4	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	0.5
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
Ethanol	ND	1.0	50	Ethylbenzene	ND	1.0	0.5
Ethyl tert-butyl ether (ETBE)	ND	1.0	0.5	Hexachlorobutadiene	ND	1.0	0.5
2-Hexanone	ND	1.0	0.5	Iodomethane (Methyl iodide)	ND	1.0	5.0
Isopropylbenzene	ND	1.0	0.5	4-Isopropyl toluene	ND	1.0	0.5
Methanol	ND	1.0	500	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	ND	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 Acetate	ND	1.0	5.0	Vinyl Chloride	ND	1.0	0.5
Xylenes	ND	1.0	0.5				
		Sur	rogate Re	ecoveries (%)			
%SS1:	111		%SS2:	100		•	
%SS3:	99.3	3	•				
Comments				Transport of the second of the			

Comments

h) lighter than water immiscible sheen/product is present; i) liquid sample that contains greater than ~2 vol. % sediment; j) sample diluted due to high organic content.



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

ND means not detected above the reporting limit; N/A means analyte not applicable to this analysis.

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

QC SUMMARY REPORT FOR SW8021B/8015Cm

Matrix: W

WorkOrder: 0311385

EPA Method: SW8	3021B/8015Cm E	Extraction: SW5030B			BatchID: 9484			Spiked Sample ID: 0311378-003B			
	Sample	Spiked	MS*	MSD*	MS-MSD LCS		LCSD	LCS-LCSD	Acceptanc	e Criteria (%)	
	μg/L	μg/L	% Rec.	% Rec.	% RPD	% Rec.	% Rec.	% RPD	Low	High	
TPH(btex) [£]	ND	60	107	106	0.770	95.6	98.9	3.33	70	130	
MTBE	ND	10	93.3	95.2	2.01	87.2	88.5	1.50	70	130	
Benzene	ND	10	100	103	2.10	98.7	102	2.76	70	130	
Toluene	ND	10	96.3	98.1	1.85	101	104	2.70	70	130	
Ethylbenzene	ND	10	104	106	1.59	102	105	2.95	70	130	
Xylenes	ND	30	100	100	0	103	107	3.17	70	130	
%SS:	98.6	100	99.7	101	0.846	108	109	0.214	70	130	

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

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 and / or MSD spike recoveries may not be near 100% or the RPDs near 0% if; a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) if that specific sample matrix interferes with spike recovery.

[£] TPH(blex) = 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 = 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.

WorkOrder: 0311385

1,2-Dichloroethane (1,2-DCA)

1,1-Dichloroethene

Ethanol

Methanol

Toluene

%SS1:

%SS2:

%SS3:

NONE

Trichloroethene

Diisopropyl ether (DIPE)

Ethyl tert-butyl ether (ETBE)

Methyl-t-butyl ether (MTBE)

ND

ND

ND

ND

ND

ND

ND

ND

ND

107

101

98.0

10

10

10

500

10

2500

10

10

10

100

100

100

95.5

74.5

92

82.4

83.3

102

86.1

100

85.1

103

100

96.4

OC SUMMARY REPORT FOR SW8260B

Matrix: W

EPA Method: SW8260B Extraction: SW5030B BatchID: 9472 Spiked Sample ID: 0311376-002C Sample Spiked MS* MSD* MS-MSD1 LCS LCSD LCS-LCSD Acceptance Criteria (%) % RPD % Rec. % Rec. % Rec. % RPD Low High μg/L µg/L % Rec. tert-Amyl methyl ether (TAME) ND 10 82.8 83.9 1.27 87.6 82.1 6.54 70 130 Benzene ND 10 88.1 89.8 1.86 92.4 89.9 2.66 70 130 t-Butyl alcohol (TBA) ND 50 84.9 83 2.19 88.3 81.1 8.47 70 130 Chlorobenzene ND 10 94.9 97.1 2.24 99.6 95.3 4.37 70 130 1,2-Dibromoethane (EDB) ND 10 107 70 105 1.31 111 106 5.36 130

98

77.7

94.5

98.6

85.4

102

89

101

86.7

105

100

95.9

2.49

4.19

2.69

17.9

2.46

0

3.26

0.943

1.81

1.98

0

0.568

101

79.4

98

82.6

89

98.8

90.4

104

89.5

99.6

100

95.7

96.8

78.9

92.1

88.1

83.8

98.6

84.2

99.4

86.7

99.6

100

97.8

3.81

0.712

6.24

6.40

5.96

0.191

7.11

4.88

3.22

0

0

70

70

70

70

70

70

70

70

70

70

70

130

130

130

130

130

130

130

130

130

130

130

130

70 2.25 All target compounds in the Method Blank of this extraction batch were ND less than the method RL with the following exceptions:

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 and / or MSD spike recoveries may not be near 100% or the RPDs near 0% if: a) the sample is inhomogenous AND contains significant concentrations of analyte relative to the amount spiked, or b) if that specific sample matrix interferes with 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 ianalyte content.

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

Q.	P & D ENVIR A Division of Paul 4020 Panam Oakland, C. (510) 658	H, King, Inc. ia Court A 94611	ΓAL	C	CHAL	N OF (CUSTOE)Y F	3F)	CO	RD:	* 3°	B (03113	SS GE <u>l</u> OF	
!!	PROJECT NUMBER: 0055 SAMPLED BY: (PRI WILLENUMBER	NTED AND	signat back	TOrm	er H	aber Oil		NUMBER OF CONTAINERS	7 DILL YSISTES				PRESED	/	GE OF	
+ ++++	MWI MWZ MW3 MW5 MW6 MW7 MW8	11/25/03		wher				5	XXXXXXX	X	I I	CONDITI PACE A ORINAT	DN. SERVIT	APPR	OPRIATE ALS OTHER	
	RELINQUISHED BY: RELINQUISHED BY: RELINQUISHED BY:	2000 (SIGNATURE #22)	<u>(</u>)	DATE DATE DATE	TIME 10:03 TIME	RECEIVED BY: NULL VOL RECEIVED FOR (SIGNATURE)	(SIGNATURE) # 280. (SIGNATURE) 1/25 R LABORATORY	8:30ph 8Y:	TOTAL TOTAL LAB	THIS SHEP NO. OF THIS SHEP BORATO S	CONTAINE MENT) ORY C IO R SAMPLE	ONTAC ONTAC MACHED:	O M CT: LAB MC (9 LYSIS R ()YE		M Anal PHONE NUT 18 - 16-2 HEET	MBER:

CHAIN-OF-CUSTODY RECORD

Page I of

110 Second Avenue South, #D7 Pacheco, CA 94553-5560 (925) 798-1620

WorkOrder: 0311385

Client:

P & D Environmental 4020 Panama Court Oakland, CA 94611-4931 TEL:

(510) 658-6916

FAX:

(510) 834-0772

ProjectNo:

#0055; Former Haber Oil

PO:

Date Received:

11/26/03

Date Printed:

11/26/03

				1		Requested Tests					
Sample ID	ClientSampID	Matrix	Collection Date	Hold	8260B+7OXY_W	G-MBTEX_W					
0311385-001	MW 1	Water	11/25/03	ľ 🗀 !	B :						
0311385-002	MW 2	Water	11/25/03		В	Α					
0311385-003	MW 3	Water	11/25/03		В	Α	,				
0311385-004	MW 4	Water	11/25/03		В	A	The state of the s				
0311385-005	MW 5	Water	11/25/03		В	A					
311385-006	MW 6	Water	11/25/03		В	Α					
311385-007	MW 7	Water	11/25/03		B	Α					
311385-008	MW B	Water	11/25/03		В	Α	100 M.C. VIII				

Prepared	by:	Melissa	Valle
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Comments:

NOTE: Samples are discarded 60 days after results are reported unless other arrangements are made. Hazardous samples will be returned to client or disposed of at client expense.