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Alameda County Environmental Health

Groundwater Monitoring Report January 1 through March 31, 2009 Former Hot Mix Asphalt Plant Area (AOC #1) Hanson Aggregates Radum Facility 3000 Busch Road, Pleasanton, California (ACEH Case #RO0002941 and Geotracker Global ID #SLT19719376)

May 11, 2009 001-09567-07

Prepared for: Lehigh Hanson West Region 12667 Alcosta Boulevard, Suite 400 San Ramon, California 94583

Prepared by LFR Inc. 1900 Powell Street, 12th Floor Emeryville, California 94608



May 11, 2009

Mr. Jerry Wickham Alameda County Health Care Services Environmental Health Services 1131 Harbor Bay Parkway, Suite 250 Alameda, California 94502-6577

Subject: Groundwater Monitoring Report, January 1 through March 31, 2009, Former

Hot Mix Asphalt Plant Area (AOC #1), Hanson Aggregates Radum Facility, 3000 Busch Road, Pleasanton, California (ACEH Case #RO0002941 and

Geotracker Global ID # SLT19719376)

Dear Mr. Wickham:

The enclosed Groundwater Monitoring Report was prepared by LFR Inc. (LFR) on behalf of Lehigh Hanson West Region for the former hot mix asphalt plant area (located within Area of Concern [AOC] #1) of the Hanson Aggregates Radum Facility, located at 3000 Busch Road, Pleasanton, California ("the Site"). This report presents and discusses the results of the fourth of four planned quarterly groundwater monitoring events conducted at the Site.

This event and the previous three quarterly groundwater monitoring events were conducted in accordance with the February 28, 2008 work plan approved by Alameda County Environmental Health in its technical comment letter dated March 31, 2008. Results from the groundwater monitoring confirm that groundwater beneath the Site has not been affected by total petroleum hydrocarbons previously detected in soil in limited areas of the Site. Based on the results of quarterly groundwater monitoring, and in accordance with the work plan, LFR recommends that groundwater monitoring be discontinued and that the 10 groundwater monitoring wells be properly abandoned.

I declare, under penalty of perjury, that the information and/or recommendations contained in the attached Report are true and correct to the best of my knowledge.

If you have any questions or comments concerning this report, please call me at (925) 244-6584 or Ron Goloubow of LFR at (510) 652-4500.

Sincerely,

Lee W. Cover

Environmental Manager

Lee W. an

Hanson Aggregates Northern California

Attachment

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CERTIFICATIONS

LFR Inc. has prepared this Groundwater Monitoring Report on behalf of Lehigh Hanson West Region in a manner consistent with the level of care and skill ordinarily exercised by professional geologists and environmental scientists. This report was prepared under the technical direction of the undersigned California Professional Geologist.





May 11, 2009

Katrin Schliewen, P.G.

Senior Hydrogeologist

California Professional Geologist (7808)

Date

Ron Goloubow

Senior Associate Geologist

May 11, 2009
Date

EXECUTIVE SUMMARY

This Groundwater Monitoring Report for the period January 1 through March 31, 2009 ("the reporting period") presents the results of the fourth quarterly groundwater monitoring event conducted by LFR Inc. (LFR) in the former hot mix asphalt plant area of the Hanson Aggregates Radum Facility located in Pleasanton, California ("the Site").

The groundwater monitoring event that was completed during this reporting period represents the fourth of four required periodic groundwater monitoring events for the Site. The first, second, and third groundwater monitoring events were conducted during June 2008, September 2008, and January 2009, respectively. The fourth groundwater monitoring event consisted of measuring depth to groundwater and purging and sampling wells MW-1 through MW-10 and was conducted on March 16 and 17, 2009. Well MW-4 could not be sampled due to an insufficient amount of water in the well.

Equipotential contours drawn based on groundwater elevations indicate that the local groundwater flow direction is toward the northwest, with a horizontal groundwater gradient of approximately 0.018 foot per foot. Analytical results indicate that none of the analyzed constituents were detected above laboratory reporting limits in any of the groundwater samples collected during this groundwater monitoring event. The analytical results from the current sampling event are consistent with results from previous groundwater monitoring events.

Based on the results of the four quarterly groundwater monitoring events conducted at the Site, and in accordance with the February 28, 2008 work plan approved by Alameda County Environmental Health, LFR recommends that periodic groundwater monitoring be discontinued and that the 10 groundwater monitoring wells be properly abandoned.

1.0 INTRODUCTION

This Groundwater Monitoring Report presents the results of groundwater monitoring activities conducted for the period from January 1 through March 31, 2009 ("the reporting period") by LFR Inc. (LFR) on behalf of Lehigh Hanson West Region ("Hanson") in the former hot mix asphalt plant area of the Hanson Aggregates Radum Facility located at 3000 Busch Road, Pleasanton, California ("the Site"; Figures 1 and 2). The Site is located within Area of Concern #1 (AOC #1). Quarterly groundwater monitoring was conducted in the 10 groundwater monitoring wells installed at the Site during 2007 and 2008 (wells MW-1 through MW-10) to monitor groundwater quality and groundwater flow direction and gradient for approximately one year.

Quarterly groundwater monitoring was initiated at AOC #1 in accordance with the scope of work described in the "Work Plan for Additional Well Installations and Quarterly Groundwater Monitoring and Reporting in the Former Hot Mix Asphalt Plant Area (AOC #1) of the Hanson Aggregates Radum Facility, 3000 Busch Road, Pleasanton, California, SLIC Case #RO0002941 and Geotracker ID SLT19719376" ("the Work Plan"), which was submitted to Alameda County Environmental Health (ACEH) on February 28, 2008. ACEH approved the Work Plan on March 31, 2008. In its approval letter, ACEH requested that sampling for dissolved metals be conducted during the second quarterly groundwater monitoring event instead of the first as proposed in the Work Plan, to allow additional time to pass between installing the last three groundwater monitoring wells and sampling for dissolved metals.

With this reporting period, LFR has completed the four planned approximately quarterly groundwater monitoring events on behalf of Hanson, which were conducted during June and September 2008, and January and March 2009. The March 2009 groundwater monitoring event that was completed during this reporting period represents the fourth of the four required quarterly events for the Site. Results from this event and the three previous reporting periods are presented in this report. Based on the analytical results for samples collected during the four groundwater monitoring events, no additional quarterly groundwater monitoring events are planned for the Site.

2.0 METHODOLOGY

2.1 Quarterly Groundwater Monitoring

This quarterly groundwater monitoring event was conducted on March 16 and 17, 2009, and consisted of measuring depth to groundwater and of purging and sampling groundwater monitoring wells MW-1 through MW-10 (Table 1 and Figure 3). The methodology of the groundwater monitoring event is described in this section, and results are presented and discussed in Section 3.0.

2.1.1 Groundwater Elevation Monitoring

Groundwater elevation monitoring was conducted at AOC #1 on March 16, 2009. The depth to groundwater was measured prior to purging and sampling, using a Solinst water level indicator, and relative to the top of casing (TOC). Depth-to-groundwater measurements were recorded on field sheets, copies of which are included in Appendix A. Groundwater elevations were calculated by subtracting the depth-to-groundwater measurement from the TOC elevation (Table 3).

2.1.2 Groundwater Well Purging and Sampling

Wells MW-1 through MW-10 (except for well MW-4) were purged and sampled on March 16 and 17, 2009, using single-use, disposable bailers. Consistent with previous sampling events, well MW-4 did not contain a sufficient amount of groundwater for purging and sampling. The Work Plan proposed that a low-flow purging and sampling technique would be used during the quarterly groundwater monitoring events. While implementing this purging method during the first quarterly monitoring event, it was discovered that the yield for many of the wells is not sufficient to sustain a consistent water level during low-flow sampling; however, the wells could produce three casing volumes. Therefore, three casing volumes of groundwater were purged from each well using disposable bailers before the groundwater sample was collected. This method of purging and sample collection was employed during each of the four quarterly groundwater sampling events.

Depth-to-groundwater and general water-quality parameters were monitored during purging, and the parameters were recorded on field sheets, copies of which are included in Appendix A. The wells were considered sufficiently purged after at least three casing volumes were removed from each well and general water-quality parameters stabilized. Groundwater samples were collected after purging was completed.

Groundwater samples were collected in clean, laboratory-provided sample containers, properly labeled, and stored in an ice-chilled cooler for transport to the analytical laboratory under chain-of-custody protocol. One field duplicate sample was collected from well MW-8. In addition, a field blank was collected and submitted to the laboratory for quality control purposes.

2.1.3 Quarterly Monitoring Laboratory Analyses

Groundwater samples were submitted to TestAmerica Laboratories, Inc., a California-certified analytical laboratory located in Pleasanton, California. All samples were analyzed for the following parameters, and in accordance with the sample matrix presented in Table 2:

- Total petroleum hydrocarbons (TPH) as diesel (TPHd) and TPH as motor oil (TPHmo) by U.S. Environmental Protection Agency (EPA) Method 8015 (after undergoing silica-gel cleanup)
- TPH as gasoline (TPHg) by EPA Method 8260
- Benzene, toluene, ethylbenzene, and total xylenes (BTEX) by EPA Method 8260
- Fuel oxygenates by EPA Method 8260
- Lead scavengers by EPA Method 8260
- Semivolatile organic compounds (SVOCs) by EPA Method 8270

3.0 RESULTS

Results from the current reporting period and from previous quarterly groundwater monitoring events are presented in this report. Historical groundwater elevation data are summarized in Table 3, and interpreted groundwater equipotential contours for the reporting period are presented on Figure 4. Historical analytical results are summarized in Table 4, and analytical results for the reporting period are presented on Figure 5. Relevant San Francisco Bay Regional Water Quality Control Board Environmental Screening Levels (ESLs) are included in the summary tables, and compounds detected at concentrations that exceeded the ESLs are highlighted in the appropriate summary tables and figures.

3.1 Groundwater Elevations

The groundwater elevation contours indicate that the groundwater flow direction beneath the Site was approximately to the northwest on March 16, 2009, with a horizontal groundwater gradient of approximately 0.018 foot per foot (Figure 4).

Groundwater elevations for this monitoring period are consistent with results from previous groundwater monitoring events. Monitoring well MW-4 has been dry since it was installed and was dry during this monitoring event. Over the four monitoring events, groundwater elevations generally have increased approximately 1 to 2 feet, with the exception of well MW-5 where the groundwater elevation has decreased by approximately 3 feet. Monitoring well MW-5 was installed with a well screen approximately 10 to 15 feet deeper than most of the other wells at the Site because of subsurface conditions encountered during drilling. The groundwater elevation from this well historically has been approximately 20 feet lower than other wells at the Site and therefore has not been used for creating the equipotential contours.

3.2 Groundwater Analytical Results

Analytical results indicate that none of the compounds analyzed for in the groundwater monitoring wells were detected above laboratory reporting limits during this

groundwater monitoring event (Table 4 and Figure 5). These results are consistent with analytical results from previous quarterly groundwater sampling events at the Site. During the quarterly groundwater monitoring events, the primary contaminants of concern (COCs) at the Site (TPHd and TPHmo) have not been detected above laboratory reporting limits in any groundwater samples collected from the existing groundwater monitoring wells. Similarly, TPHg, fuel oxygenates, and lead scavengers have not been detected in any groundwater samples collected from the groundwater monitoring wells.

As shown in Table 4, only sporadic low concentrations of non-COCs have been reported for groundwater samples collected approximately quarterly from groundwater monitoring wells at the Site. Historically, the only TPH-related compound detected above laboratory limits has been toluene detected at low concentrations (estimated below the laboratory reporting limit) in the October 2007 samples collected from wells MW-3 (0.3 microgram per liter [μ g/l]) and MW-5 (0.4 μ g/l). None of the BTEX compounds have been detected since the estimated toluene concentrations were reported for the two October 2007 samples.

Of the other compounds analyzed in groundwater samples, only one SVOC and two dissolved metals have been detected above laboratory reporting limits. A low concentration of bis(2-thylhexyl)phthalate was detected in the primary sample collected from well MW-2 (9.8 μ g/l) in September 2008, but this compound was not detected in the duplicate sample collected from the same well. Low concentrations of barium (below the ESL of 1,000 μ g/l for barium) were detected in the September 2008 samples collected from wells MW-3 (160 μ g/l), MW-8 (230 μ g/l), and MW-9 (150 μ g/l), and a concentration of copper (slightly above the ESL of 3.1 μ g/l for copper) was detected in the September 2008 sample collected from well MW-9 (5 μ g/l).

The analytical results from the reporting period, and from the previous three quarterly groundwater monitoring events, confirm that groundwater beneath the Site has not been affected by the TPH or TPH-related compounds that have been detected in soil in limited areas beneath the Site.

4.0 CONCLUSIONS AND RECOMMENDATIONS

4.1 Conclusions

The primary COCs detected in soil and groundwater during previous investigations are TPHd, and to a lesser extent TPHmo. Other compounds, including TPHg, volatile organic compounds (VOCs), BTEX, SVOCs, and metals, have been detected only sporadically in isolated soil or groundwater samples. Fuel oxygenates, lead scavengers, pesticides, and polychlorinated biphenyls (PCBs) have not been detected in any samples collected. Only concentrations detected above the May 2008 ESLs are considered significant. Based on results from previous investigations, LFR has concluded that

TPH-affected soil is limited in extent both vertically and laterally and that groundwater has not significantly been affected by previous operations conducted at the Site (LFR 2007). The four planned quarterly groundwater monitoring events consisting of measuring depth to groundwater and collecting groundwater samples from 10 existing groundwater monitoring wells installed at the Site have been concluded.

The results from the four completed quarterly groundwater monitoring and sampling events confirm that groundwater beneath the Site has not been significantly affected by TPH or TPH-related compounds detected during previous investigations. Groundwater elevation data indicate that the direction of groundwater flow beneath the Site was consistently to the west-northwest, with a horizontal groundwater gradient ranging from approximately 0.015 to 0.025 foot per foot. The primary COCs (TPHd and TPHmo) have not been detected above laboratory reporting limits in groundwater samples collected from the groundwater monitoring wells during the four quarterly groundwater monitoring events. Similarly, TPHg, fuel oxygenates, and lead scavengers have not been detected in samples collected from groundwater monitoring wells. Other compounds, including toluene, bis(2-ethylhexyl)phthalate, and dissolved barium and copper, have been detected only once in up to three wells (in the case of barium). None of these low detections are considered significant because they were estimated at concentrations below laboratory reporting limits, are below the ESLs, were not detected in duplicate samples, and/or were not detected in subsequent groundwater samples.

4.2 Recommendations

Based on results of the four quarterly groundwater monitoring events completed at the Site, and in accordance with the Work Plan, LFR recommends that quarterly groundwater monitoring and reporting be discontinued and that the 10 groundwater monitoring wells be properly abandoned.

4.3 Future Activities

LFR understands that the Site may in the future undergo a property transfer to Legacy Partners ("Legacy"), similar to the property transfer agreement between Hanson and Legacy that was completed for the majority of the Radum property during 2007. It is assumed that under Legacy the Site eventually would be developed for commercial and light-industrial land use.

In accordance with ongoing discussions between Hanson and ACEH (for subsurface environmental concerns) and the Livermore-Pleasanton Fire Department (LPFD; for remaining surface features and associated materials), certain closure activities will be conducted prior to a potential land transfer to Legacy. Additional subsurface investigations are proposed to be conducted in limited areas of the Site as described in the September 15, 2008 "Work Plan for Additional Site Characterization at Selected Areas Within AOC #1" approved by ACEH on January 29, 2009. Based on the results of this planned and other previously completed subsurface investigations, affected soil

will be excavated from selected areas and confirmation sampling will be conducted. At the request of the LPFD, Hanson will submit a closure plan to describe planned closure activities for surface features, including the former paving oil containment structure and its contents, the former truck scale, and other features planned for removal and/or disposal by Hanson.

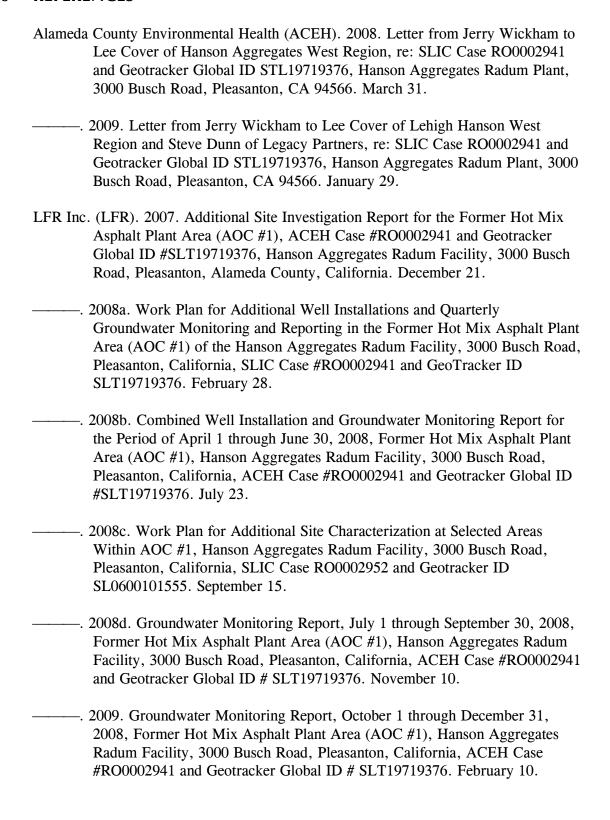
5.0 LIMITATIONS

The opinions and recommendations presented in this report are based upon the scope of services, information obtained through the performance of the services, and the schedule as agreed upon by LFR and the party for whom this report was originally prepared. This report is an instrument of professional service and was prepared in accordance with the generally accepted standards and level of skill and care under similar conditions and circumstances established by the environmental consulting industry. No representation, warranty, or guarantee, expressed or implied, is intended or given. To the extent that LFR relied upon any information prepared by other parties not under contract to LFR, LFR makes no representation as to the accuracy or completeness of such information. This report is expressly for the sole and exclusive use of the party for whom this report was originally prepared for a particular purpose. Only the party for whom this report was originally prepared and/or other specifically named parties have the right to make use of and rely upon this report. Reuse of this report or any portion thereof for other than its intended purpose, or if modified, or if used by third parties, shall be at the user's sole risk.

Results of any investigations or testing and any findings presented in this report apply solely to conditions existing at the time when LFR's investigative work was performed. It must be recognized that any such investigative or testing activities are inherently limited and do not represent a conclusive or complete characterization. Conditions in other parts of the Site may vary from those at the locations where data were collected. LFR's ability to interpret investigation results is related to the availability of the data and the extent of the investigation activities. As such, 100 percent confidence in environmental investigation conclusions cannot reasonably be achieved.

LFR, therefore, does not provide any guarantees, certifications, or warranties regarding any conclusions regarding environmental contamination of any such property. Furthermore, nothing contained in this document shall relieve any other party of its responsibility to abide by contract documents and applicable laws, codes, regulations, or standards.

6.0 REFERENCES



Regional Water Quality Control Board, San Francisco Bay Region. 2008. Screening for Environmental Concerns at Sites with Contaminated Soil and Groundwater (Interim Final – May 2008); Environmental Screening Levels ("ESLs"). Technical Document. May.

Table 1
Groundwater Monitoring Well Construction Details
Area of Concern # 1 / Former Hot Mix Asphalt Plant Area
Hanson Radum Facility, 3000 Busch Road, Pleasanton, California

Monitoring Well ID	Installation Date	Boring Hole Diameter (inches)	Casing Diameter (inches)	Approximate Total Well Depth (feet bgs)	Approximate Screened Interval (feet bgs)	Top of Casing Elevation ¹ (feet msl)
MW-1	10/3/07	8.0	2.0	60	45 - 60	374.67
MW-2	10/2/07	8.0	2.0	60	45 - 60	376.33
MW-3	10/4/07	8.0	2.0	60	45 - 60	374.95
MW-4	10/5/07	8.0	2.0	48	43 - 48	372.94
MW-5	10/9/07	8.0	2.0	74	69 - 74	374.35
MW-6	10/10/07	8.0	2.0	55	45 - 55	375.03
MW-7	10/1/07	8.0	2.0	65	50 - 65	377.68
MW-8	6/9/08	8.0	2.0	61	51 - 61	378.60
MW-9	6/10/08	8.0	2.0	52	42 - 52	375.75
MW-10	6/11/08	8.0	2.0	54	44 - 54	375.62

Notes:

ID = identification; monitoring well identification number

feet bgs = feet below ground surface

feet msl = feet relative to mean sea level

¹ Top of casing elevation is approximately 3.0 feet above ground surface; top of casing elevation and land survey conducted by Kier & Wright Civil Engineers & Surveyors, Inc.

Table 2
Quarterly Groundwater Monitoring Sample Matrix
Area of Concern #1 / Former Hot Mix Asphalt Plant Area
Hanson Radum Facility, 3000 Busch Road, Pleasanton, California

Well ID	Date Installed	Approximate Top (feet bgs)	Screen Interval Bottom (feet bgs)	TPHd / TPHmo 8015	TPHg 8260	BTEX 8260	Fuel Ox 8260	Lead Scav 8260	SVOCs 8270	Dissolved Metals 6010B
Groundwater Monit	oring Wells									
MW-1	10/3/2007	45	60	X	X	X	X	X	X	-
MW-2	10/2/2007	45	60	X	X	X	X	X	X	-
MW-3	10/4/2007	45	60	X	X	X	X	X	X	once 1
MW-4	10/5/2007	43	48	X	X	X	X	X	X	-
MW-5	10/9/2007	69	74	X	X	X	X	X	X	-
MW-6	10/10/2007	45	55	X	X	X	X	X	X	-
MW-7	10/1/2007	50	65	X	X	X	X	X	X	-
MW-8	6/9/2008	51	61	X	X	X	X	X	X	once 1
MW-9	6/10/2008	42	52	X	X	X	X	X	X	once 1
MW-10	6/11/2008	44	54	X	X	X	X	X	X	-
Quality Assurance and Quality Control Samples ²										
Field Blank	na	na	na	X	X	X	X	X	X	-
Trip Blank	na	na	na	-	X	X	X	X	-	-

Notes:

feet bgs = feet below ground surface

na = not applicable

EPA = U.S. Environmental Protection Agency

TPHd = total petroleum hydrocarbons as diesel by EPA Method 8015 (with silica-gel cleanup)

TPHmo = total petroleum hydrocarbons as motor oil by EPA Method 8015 (with silica-gel cleanup)

TPHg = total petroleum hydrocarbons as gasoline by EPA Method 8260

BTEX = benzene, toluene, ethylbenzene, and total xylenes by EPA Method 8260

Fuel Ox = fuel oxygenates by EPA Method 8260

Lead Scav = lead scavengers by EPA Method 8260

SVOCs = semivolatile organic compounds by EPA Method 8270

Dissolved Metals = CAM 17 list of dissolved metals (laboratory filtered samples) by EPA Method 6010B

[&]quot;x" = to be analyzed quarterly for four consecutive quarters

[&]quot;-" = not analyzed

¹ Samples for dissolved metals were collected only once, during the second groundwater monitoring event conducted on September 16, 2008.

² One field blank (FB) sample was collected during each quarterly monitoring event, and one trip blank (TB) sample was collected for every cooler of samples transported to the laboratory during every quarterly monitoring event.

Table 3
Groundwater Elevations
Area of Concern # 1 / Former Hot Mix Asphalt Plant Area
Hanson Radum Facility, 3000 Busch Road, Pleasanton, California

Monitoring Well ID	Top of Casing Elevation ¹ (feet msl)	Measurement Date	Depth to Groundwater Measured (feet TOC)	Groundwater Elevation (feet msl)
MW-1	374.67	10/22/07	57.22	317.45
		6/16/08	57.35	317.32
		9/15/08	57.59	317.08
		1/9/09	57.79	316.88
		3/16/09	55.92	318.75
MW-2	376.33	10/22/07	55.24	321.09
		6/16/08	55.39	320.94
		9/15/08	55.73	320.60
		1/9/09	56.05	320.28
		3/16/09	54.56	321.77
MW-3	374.95	10/22/07	54.32	320.63
		6/16/08	54.53	320.42
		9/15/08	54.74	320.21
		1/9/09	55.00	319.95
		3/16/09	53.64	321.31
MW-4	372.94	10/22/07	47.37	325.57
		6/16/08	48.77	324.17
		9/15/08	48.71	324.23
		1/9/09	48.85	DRY
		3/16/09	48.54	DRY
MW-5	374.35	10/22/07	68.40	305.95
		6/16/08	70.16	304.19
		9/15/08	70.16	304.19
		1/9/09	75.04	299.31
		3/16/09	71.36	302.99
MW-6	375.03	10/22/07	49.19	325.84
		6/16/08	49.34	325.69
		9/15/08	49.49	325.54
		1/9/09	49.67	325.36
		3/16/09	48.89	326.14
MW-7	377.68	10/22/07	57.04	320.64
		6/16/08	57.21	320.47
		9/15/08	57.79	319.89
		1/9/09	58.14	319.54
		3/16/09	56.54	321.14
MW-8	378.60	10/22/07		NM
		6/16/08	55.73	322.87
		9/15/08	55.99	322.61
		1/9/09	56.22	322.38
		3/16/09	54.61	323.99

Table 3
Groundwater Elevations
Area of Concern # 1 / Former Hot Mix Asphalt Plant Area
Hanson Radum Facility, 3000 Busch Road, Pleasanton, California

Monitoring Well ID	Top of Casing Elevation ¹ (feet msl)	Measurement Date	Depth to Groundwater Measured (feet TOC)	Groundwater Elevation (feet msl)
MW-9	375.75	10/22/07	 51 40	NM
		6/16/08 9/15/08	51.48 51.71	324.27 324.04
		1/9/09	51.88	323.87
		3/16/09	50.45	325.30
MW-10	375.62	10/22/07		NM
		6/16/08	51.38	324.24
		9/15/08	51.58	324.04
		1/9/09	51.64	323.98
		3/16/09	50.35	325.27

Notes:

ID = identification; monitoring well identification number

feet msl = feet relative to mean sea level

feet TOC = feet below top of casing

NM = not measured because well was not installed until June 2008

¹ Top of casing elevation is approximately 3.0 feet above ground surface; top of casing elevation and land survey conducted by Kier & Wright Civil Engineers & Surveyors, Inc.

Table 4
Groundwater Analytical Results
Area of Concern # 1 / Former Hot Mix Asphalt Plant Area
Hanson Radum Facility, 3000 Busch Road, Pleasanton, California

Monitoring Well		Total P	etroleum Hydroc	arbons					Dissolved
ID	Date	TPHd* (ug/l)	TPHmo* (ug/l)	TPHg (ug/l)	BTEX (ug/l)	Fuel Ox (ug/l)	Lead Scav (ug/l)	SVOCs (ug/l)	Metals (ug/l)
MW-1	10/22/07	< 50	< 300	< 50	ND	ND	ND	-	-
	6/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	9/15/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	1/13/09	< 50	< 300	< 50	ND	ND	ND	ND	-
	3/16/09	< 50	< 300	< 50	ND	ND	ND	ND	-
MW-2	10/22/07	< 50	< 300	< 50	ND	ND	ND	_	_
	6/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	9/15/08	< 50/ < 50	<300/<300	< 50/ < 50	ND / ND	ND / ND	ND / ND	ND ³ /ND	-
	1/9/09	< 50	< 300	< 50	ND	ND	ND	ND	-
	3/17/09	< 50	< 300	< 50	ND	ND	ND	ND	-
MW-3	10/22/07	< 50/ < 50	<300/<300	< 50/ < 50	$0.3J^{1} / 0.3J^{1}$	ND/ND	ND / ND	-	-
	6/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	9/16/08	< 50	< 300	< 50	ND	ND	ND	ND	ND ⁴
	1/13/09	< 50	< 300	< 50	ND	ND	ND	ND	-
	3/16/09	< 50	< 300	< 50	ND	ND	ND	ND	-
MW-4	10/22/07	-	-	-	-	-	-	-	-
(dry)	6/16/08	-	-	-	_	-	-	-	-
. •	9/15/08	-	-	-	_	-	-	-	-
	1/9/09	-	-	-	-	-	-	-	-
	3/17/09	-	-	-	-	-	-	-	-
MW-5	10/22/07	< 50	< 300	< 50	$0.4J^2$	ND	ND	-	-
	6/16/08	< 50/ < 50	<300/<300	< 50/ < 50	ND / ND	ND / ND	ND / ND	ND / ND	-
	9/15/08	-	-	-	_	-	-	-	-
	1/9/09	< 50	< 300	< 50	ND	ND	ND	ND	-
	3/17/09	< 50	< 300	< 50	ND	ND	ND	ND	-
MW-6	10/22/07	< 50	< 300	< 50	ND	ND	ND	-	-
	6/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	9/15/08	< 50	< 300	< 50	ND	ND	ND	ND	-

Table 4
Groundwater Analytical Results
Area of Concern # 1 / Former Hot Mix Asphalt Plant Area
Hanson Radum Facility, 3000 Busch Road, Pleasanton, California

Monitoring Well		Total P	etroleum Hydroc	arbons					Dissolved
ID	Date	TPHd* (ug/l)	TPHmo* (ug/l)	TPHg (ug/l)	BTEX (ug/l)	Fuel Ox (ug/l)	Lead Scav (ug/l)	SVOCs (ug/l)	Metals (ug/l)
MW-6	1/13/09	< 50/ < 50	<300/<300	< 50/ < 50	ND / ND	ND / ND	ND / ND	ND / ND	-
	3/17/09	< 50	< 300	< 50	ND	ND	ND	ND	-
MW-7	10/22/07	< 50	< 300	< 50	ND	ND	ND	-	-
	6/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	9/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	1/13/09	< 50	< 300	< 50	ND	ND	ND	ND	-
	3/17/09	< 50	< 300	< 50	ND	ND	ND	ND	-
MW-8	6/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	9/16/08	< 50	< 300	< 50	ND	ND	ND	ND	ND ⁴
	1/12/09	< 50	< 300	< 50	ND	ND	ND	ND	-
	3/17/09	< 50/ < 50	<300/<300	< 50/ < 50	ND / ND	ND / ND	ND / ND	ND / ND	-
MW-9	6/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	9/16/08	< 50	< 300	< 50	ND	ND	ND	ND	ND 4,5
	1/9/09	< 50	< 300	< 50	ND	ND	ND	ND	-
	3/17/09	< 50	< 300	< 50	ND	ND	ND	ND	-
MW-10	6/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	9/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	1/13/09	< 50	< 300	< 50	ND	ND	ND	ND	-
	3/17/09	< 50	< 300	< 50	ND	ND	ND	ND	-
Field Blank	6/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	9/16/08	< 50	< 300	< 50	ND	ND	ND	ND	-
	1/13/09	< 50	< 300	< 50	ND	ND	ND	ND	-
	3/17/09	< 50	< 300	< 50	ND	ND	ND	ND	-
Trip Blank	6/16/08	-	< 300	< 50	ND	ND	ND	-	_
r	9/16/08	-	< 300	< 50	ND	ND	ND	-	-
ESL	Ls (groundwater)	100	100	100	various	various	various	various	various

Table 4

Groundwater Analytical Results

Area of Concern # 1 / Former Hot Mix Asphalt Plant Area Hanson Radum Facility, 3000 Busch Road, Pleasanton, California

Monitoring Well		Total Petroleum Hydrocarbons							Dissolved
ID	Date	TPHd*	TPHmo*	TPHg	BTEX	Fuel Ox	Lead Scav	SVOCs	Metals
		(ug/l)	(ug/l)	(ug/l)	(ug/l)	(ug/l)	(ug/l)	(ug/l)	(ug/l)

Notes:

ID = identification; monitoring well identification number

ug/l = micrograms per liter

ND = not detected; no compounds were detected above their respective laboratory reporting limits

J = reported concentration is estimated below the laboratory reporting limit

"-" = sample not analyzed

"<" = not detected above the laboratory report given

TPHd = total petroleum hydrocarbons as diesel

TPHmo = total petroleum hydrocarbons as motor oil

TPHg = total petroleum hydrocarbons as gasoline

BTEX = benzene, toluene, ethylbenzene, and total xylenes

Fuel Ox = fuel oxygenates

Lead Scav = lead scavengers

SVOCs = semivolatile organic compounds

ESLs = Environmental Screening Levels by San Francisco Bay Regional Water Quality Control Board, May 2008, for Groundwater beneath Industrial/Commercial Land Use Areas where Groundwater is a Current or Potential Source of Drinking Water.

Bold font indicates that analyte detected is above the laboratory reporting limit.

Outlined values indicates that analyte was detected over the respective ESL.

^{*} Samples analyzed for TPHd and TPHmo after silica-gel cleanup except for samples collected from wells MW-2, 5, 6, 7, 8, 9, and 10 on March 17, 2009.

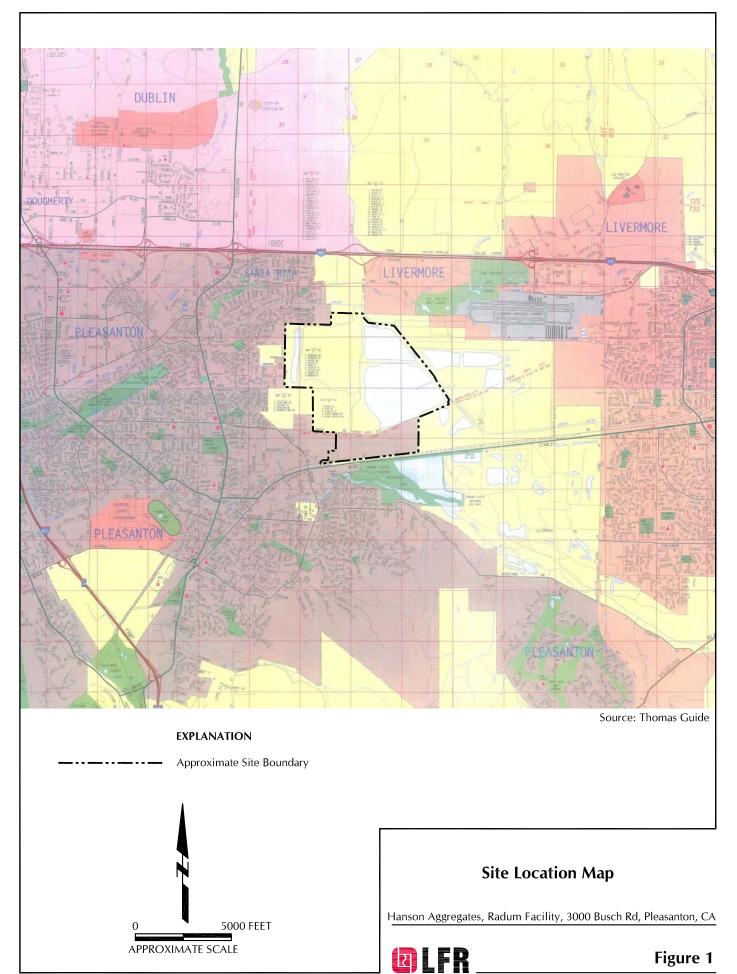
¹ Toluene was detected at a low concentration of 0.3 ug/l estimated below the laboratory reporting limit in both the primary and duplicate samples collected from well MW-3 on 10/22/07.

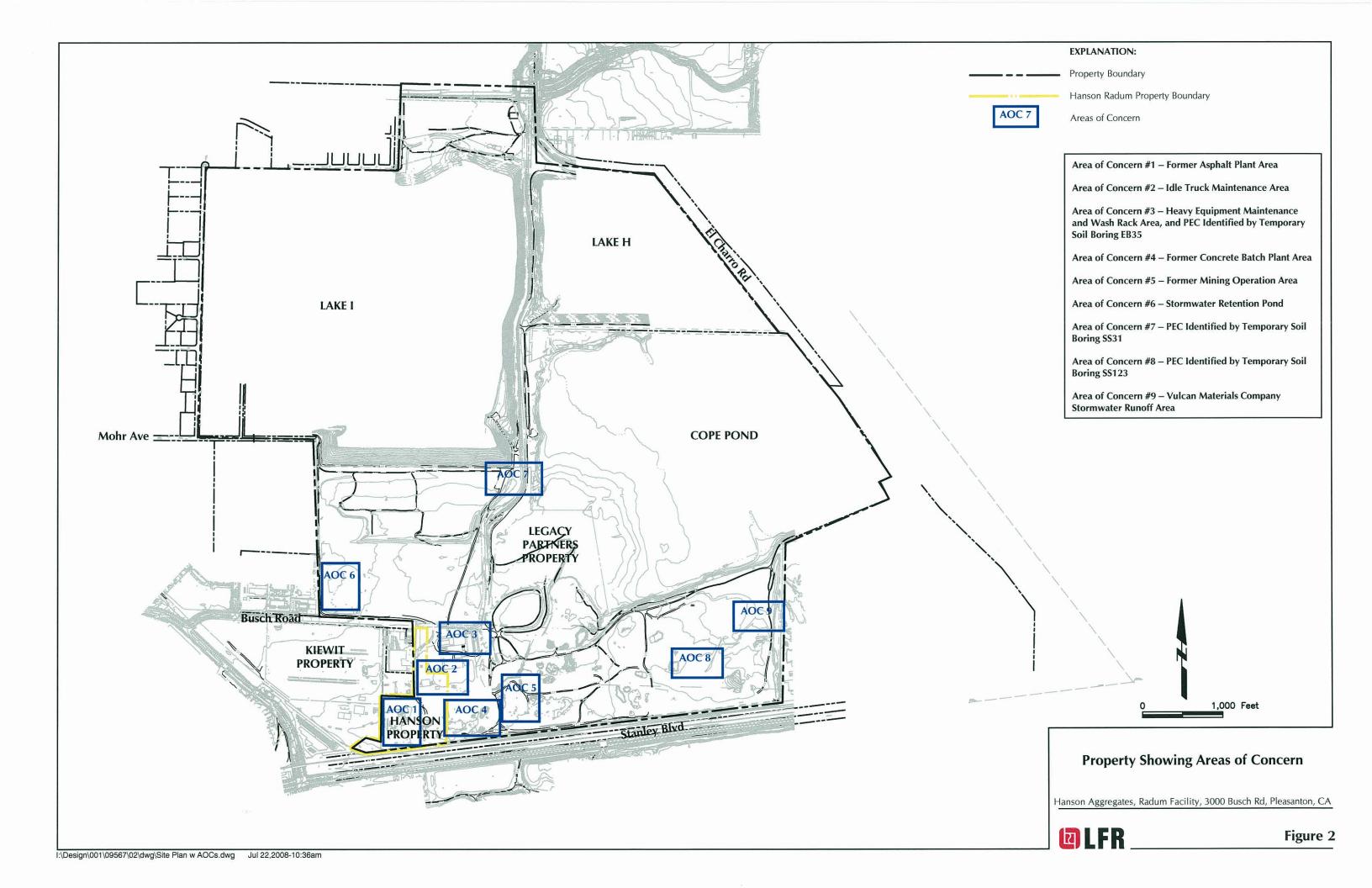
² Toluene was detected at a low concentration of 0.4 ug/l estimated below the laboratory reporting limit in the sample collected from well MW-5 on 10/22/07.

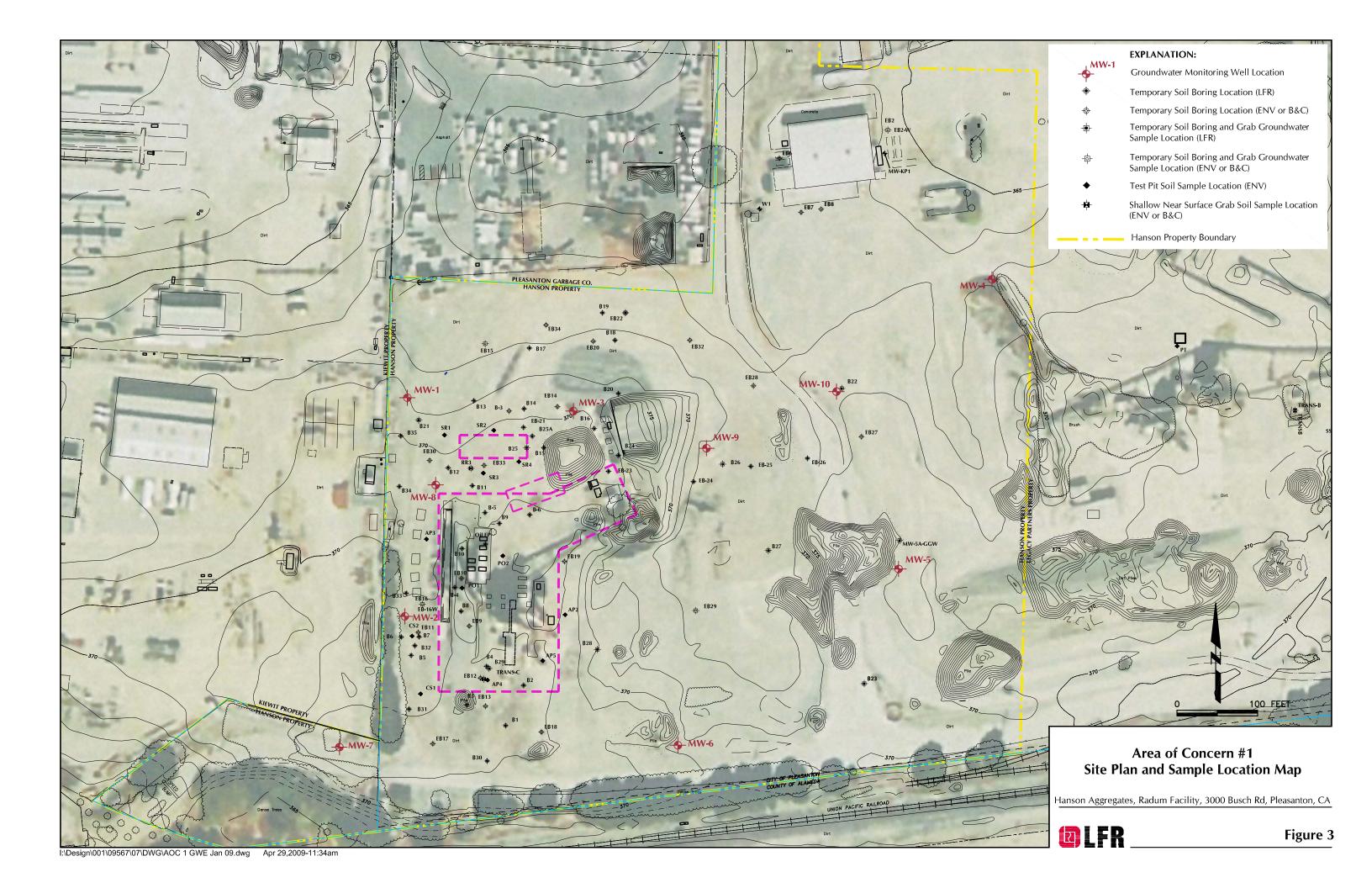
³ Bis(2-ethylhexyl)phthalate was detected at a concentration of 9.8 ug/l in the primary sample collected from MW-2 but not in the duplicate sample (ESL is 4.0 ug/l).

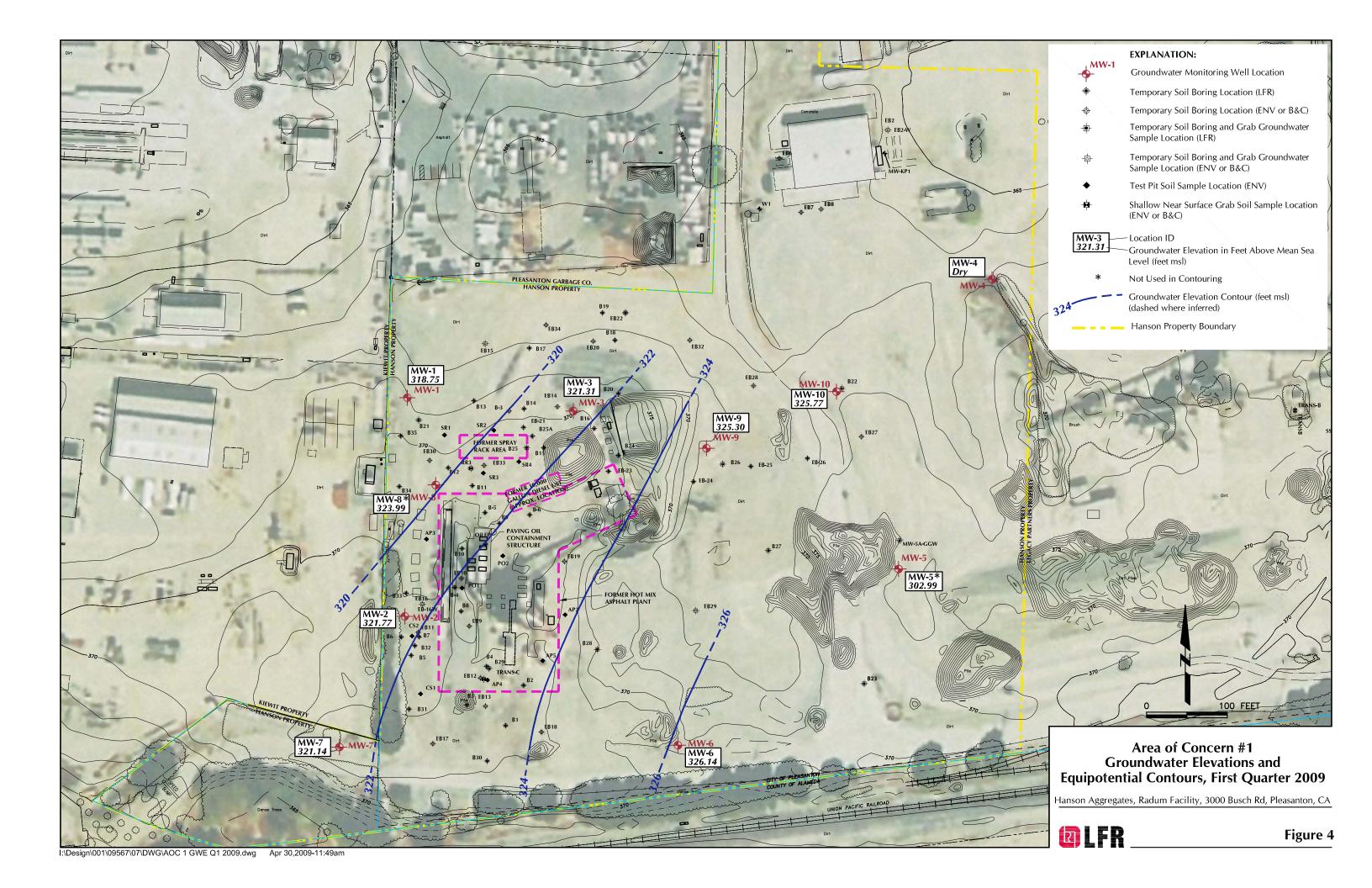
⁴ Barium was detected in the samples from wells MW-3, MW-8, and MW-9 at a concentration of 160, 230, and 150 ug/l, respectively (ESL is 1,000 ug/l).

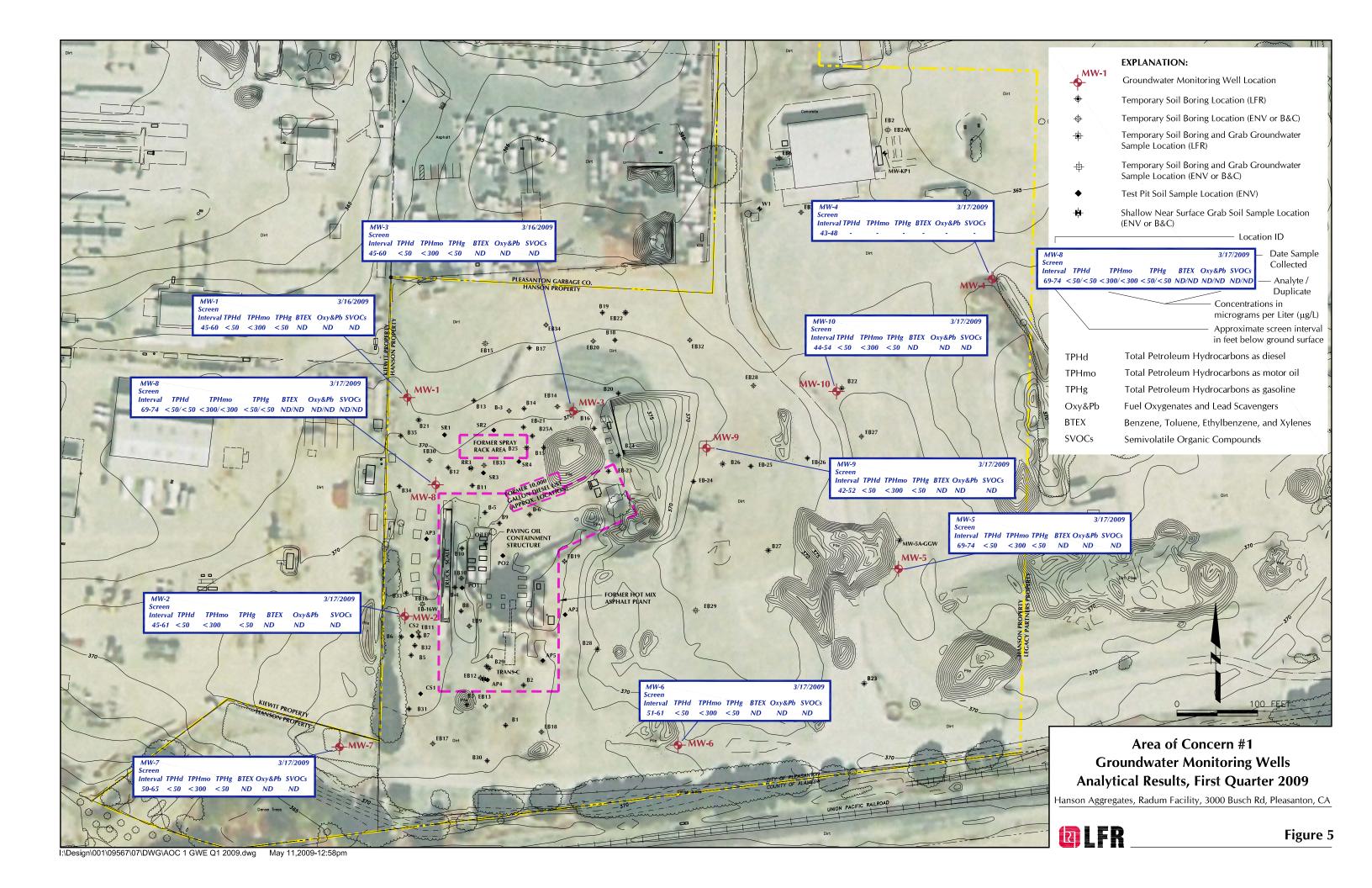
⁵ Copper was detected in sample from well MW-9 at a concentration of 5.0 ug/l (ESL is 3.1 ug/l).











APPENDIX A

Groundwater Monitoring Well Sampling Field Sheets

WATER-LEVEL MEASUREMENTS LOG

Project No. 001 - 09567 -07	Date 3/16/09	Page of	
Project Name Hanson Radum	Day: □ Sun 🙀 Mon □ Tues □ Weds		
Field Personnel 10M Cours Ashley (
General Observations Clowy			

WELL	WELL		TO WATER	WATER		ECURE?	REMARKS
NO.	ELEVATION	1	2	ELEVATION	Y	N	(UNITS = FEET)
MW-10	8:25	50.35	50.35		X		
MW-4	€:3 0	48.54	48.54		×		Hen is dry/ maddy
44-5	8:45	71.36	71.36		X		
MW-6	8:48	44.89	48.84		X		
m(n-7	8:50	56.54	56.54		*		
mu-2	8:53	54.56	54.56		人		
mw-8	8:56	54.61	54.61		#	×	Needs a JOCK
mw-1	9:00	55.92	55.92		X		
nu-3	9:03	53.64	53.64		X		
nw-9	9:05	50.46	50.45		X		
14-3		39.49	39.49		X		AOC-8
١١٠ - ١١		16.61	16.61		X		40c-8

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Project No. 001-09567-07-***	Date: March <u>/6</u> , 20	009	Page 1 of
Project Name <u>: Hanson Radum</u>	_Sampling Location: _	Hasson K	ladum
			uw-l □FB
			No.: DUP
Purge Method: ☐ Centrifugal Pump 又 Disposable Bailer 又 Hand B	ail 🛘 Submersible Pu	mp 🗷 Teflon Baile	er 🗆 Other
Purge Water Storage Container Type:	Storage Location:	On Site	
Date Purge Water Disposed:	Where Disposed:_	On Site	
Analyses Requested No. and Type	pe of Bottles Used	<u> </u>	.85-55.91
TPHg, TPHd, TPHmo First Out's		_ 6.	94 (.2)=1.388
BTEX, SVOC's, Lead Scavengers, Fuel Oxy's Lab Name: Cu rtis & Tompkin s Test	103		,
Delivery By			
Well No. MW-\ Depth of Water _5	5.41		
Well Diameter: Well Depth6	a.85		
Ż 2" (0.16 gal/feet) ☐ 5" (1.02 gal/feet) Water Column Heig			DTW 57.298
☐ 4" (0.65 gal/feet) ☐ 6" (1.47 gal/feet) Well Volume	1.11 ~1.52	80%	DIM 3 1. 2.0

Time	Inlet Depth	Depth to Water	Volume Purged (gal)	ORP	Temperature (C°)	PH (SU)	Cond (µs/cmc)	DO (mg/l)	Remarks
13:30			Sta	r+ P	415e				
B:35			1.25	-8.4	16.41	6.83	1219	2.63	
13:33			2.5	-76.7	18.28	6.96	1189	2.30	
13:38			3.75	-84.3	18.19	6.99	1184	2.74	Calibrate Sessor
13:57	:		5.0	-126.2	18.88	7.00	1203	2.37	
14:05			6.25	-117.7	18.68	7.03	1191	2.89	
14:11			7. 5	-120.4	18.53	2.01	1182	5.32	
14:20			4.75	-113	1846	7.0	1199	2.37	
14:24			10.00	-117	18.50	7.05	1188	2.45	
14:40		56.03	San	rpie					
				VRC					
					/				



Project No. 001-09567-07-***	Date: March 17, 200	Page 1 of	
Project Name <u>: Hanson Radum</u>	Sampling Location:	Hanson Radu	M
Sampler's Name: Ashley Gilleath		Sample No.: MW-2	<u> </u>
Sampling Plan By: Ron Golobow D	ated:	C.O.C. No.:	DUP
Purge Method:	ail □ Submersible Pum	np 🕱 Teflon Bailer □ Other	
Purge Water Storage Container Type:	Storage Location: C	On Site	
Date Purge Water Disposed:	Where Disposed:	On Site	-
Analyses Requested No. and Ty TPHg, TPHd, TPHmo	pe of Bottles Used	7.73	(.2) = 1.55
BTEX, SVOC's, Lead Scavengers, Fuel Oxy's Lab Name: Curtie & Fompkins TeSt Awulticar Delivery By Courier EXHand		_	
		5 80% DTW 56	.· <u>25</u>

Time	Inlet Depth	Depth to Water	Volume Purged (gal)	ORP	Temperature (C°)	PH (SU)	Cond (µs/cmc)	DO (mg/l)	Remarks
1145		Star	1 Purge						
1152			1.25	-128.8	17.93	6.94	725	3.44	
1166			a.5	-123,9			i	3.34	
1202			3.75	-121.4	17.96	685	718	3.24	
1212		54.69	Soum	ple					
	/								
				TI	20)				

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Project No. 001-09567-07-***	Date: March <u>16</u>	· · · · · · · · · · · · · · · · · · ·	Page 1 of
Project Name: Hanson Radum	Sampling Location:	TO THE PROPERTY OF THE PROPERT	
Sampler's Name: <u>Ashloy Gilrecolla</u>	Sample	No.:	D FB
Sampling Plan By: Ron Golobow	Dated:	C.O.C. No.:	DUP
Purge Method:	Bail ☐ Submersible Pump ATe	flon Bailer □ Other	
Purge Water Storage Container Type:	_ Storage Location: On Site_		
Date Purge Water Disposed:	Where Disposed: On Site		
Analyses Requested No. and T	ype of Bottles Used		
TPHg, TPHd, TPHmo			
BTEX, SVOC's, Lead Scavengers, Fuel Oxy's _			
Lab Name: Gurtis & Tompkins Test America			
Delivery By		8.78 (.2)	1.756
Well No Depth of Water _	63.64		·
Well Diameter: Well Depth 62	2. 42		
☑ 2" (0.16 gal/feet) ☐ 5" (1.02 gal/feet) Water Column He	ight <u>8.78</u>		
	1.41 ~1.5	80% DTW	5.84

Time	Inlet Depth	Depth to Water	Volume Purged (gal)	ORP	Temperature	PH (SU)	Cond (µs/cmc)	DO (mg/l)	Remarks
1325		· · · · · · · · · · · · · · · · · · ·	1.5	-7.6	63.39	6.3	. 886	367	
1337			3.0	-6.0	63.46	6.85	.772	2.26	l '
1342			4.5	8.	63.59	7.14	.768	2.98	
1346			6.0	9,6	63.95	1 -	:762	2.62	
1356			7.5	-3.3	63.63	1			
1403			9.0	5.8	63.42	7.00	,713	267	
1413			10.5	13.4	63.865	7.62	"76Z	a.33	
1424			12.0	4. H	· •	6.93	.762	1.61	
1430			13.5	-1.6	_	7.00			
1436			15.0	18.3	64.59				
1450			16.5	3.5	63.94		.756		
1458			18.0	1.2	63.58	7.01	•759	3.12	
15:05		53.69	Samo	la					
					TOC				



Project No. 001-09567-07-***	Date: March 17, 2009 Page 1 of	
Project Name <u>: Hanson Radum</u>	Sampling Location: Hapso Radum	
Sampler's Name: 10~ Collins	Sample No.: <u>MW-5</u> X FE	В
Sampling Plan By: Ron Golobow D	ed: C.O.C. No.: 🗖 DUP	
Purge Method: ☐ Centrifugal Pump ☑ Disposable Bailer ☑ Hand B	il □ Submersible Pump 🗷 Teflon Bailer □ Other	
Purge Water Storage Container Type:	Storage Location: On Site	
Date Purge Water Disposed:	Where Disposed: On Site	
Analyses Requested No. and Ty	of Bottles Used (5.90).2 =	
TPHg, TPHd, TPHmo	71.47 + 1.18	
BTEX, SVOC's, Lead Scavengers, Fuel Oxy's _	71,17 1, 1186	
Lab Name: Curtis & Tompkins TEST AMPTIZE		
Delivery By ☐ Courier And		
Well No MW - 5 Depth of Water _	7.47	
Well Diameter: Well Depth	_,	
⊠ 2" (0.16 gal/feet) □ 5" (1.02 gal/feet) Water Column Hei	, 5.90	
☐ 4" (0.65 gal/feet) ☐ 6" (1.47 gal/feet) Well Volume		

Time	Inlet Depth	Depth to Water	Volume Purged (gal)	ORP	Temperature	PH (SU)	Cond	DO (mg/l)	Remarks
8:22		. 5 t	art	Purg	e				
9:01			l.	-283	61.71	7.%	.737	4.79	
9:06			ک ک	-27.4	61.23	7.82	.741	4.77	
9:11			3	-27.9	61.63	7.78		4.36	
9:18			٧.	-23-(62.30	7.77	.725	406	
9:22			5	-71.8	61.96	7.73	-727	3 .9 0	
9:26			6	-20.5		7.75	.723		
9:30			7	-19.7	62.30	7.78	. 722	4.01	
9:35		71.48		-19.7 Scu	NR				
FB	- 02								
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Project No. 001-09567-07-***			_ Page 1 of
Project Name: Hanson Radum	Sampling Location: _	Hassot Radin	1
Sampler's Name: A Gireath		Sample No.: MW-6	□ FB
Sampling Plan By: Ron Golobow	Dated:	C.O.C. No.:	DUP
Purge Method: Centrifugal Pump Disposa	able Bailer 🗷 Hand Bail 🗆 Submersible Pur	mp 🏿 Teflon Bailer 🗆 Other	
Purge Water Storage Container Type:	•	On Site	
Date Purge Water Disposed:	Where Disposed:_	On Site	
Analyses Requested	No. and Type of Bottles Used		:
TPHg, TPHd, TPHmo			•
BTEX, SVOC's, Lead Scavengers, Fuel Oxy's _			(.2): 9.74
Lab Name: Ç urtis & Tompkin sTeS+			(.2)- 11
Delivery By	Hand		:
Well Nomw -6	Depth of Water		
Well Diameter:	Well Depth 57.94		
2" (0.16 gal/feet) ☐ 5" (1.02 gal/feet)	Water Column Height	_ - 0	3.73
☐ 4" (0.65 gal/feet) ☐ 6" (1.47 gal/feet)	Well Volume 1.44 21.5	WTQ %08	· 1.3

Time	inlet Depth	Depth to Water	Volume Purged (gal)	ORP	Temperature (C°)	PH (SU)	Cond (µs/cmc)	DO (mg/l)	Remarks
1015		Start	DOVER						
1025			1.5	-115.9	17.31	7.14	698	3,29	
1628			3	-112.9	17.25	7.11	696	3.93	
1034			4.5	-1039	17.20	7.12	686	4.48	
1037			6	-9E.	17.60	7.04	682	3.70	
1055		49.2	675	Sampl	l.				
				K 1	P_				

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Project No. 001-09567-07-***	Date: March 17, 2009 Page 1 of
Project Name <u>: Hanson Radum</u>	Sampling Location: Hasses Radum
Sampler's Name: Tom Collins	
Sampling Plan By: Ron Golobow ii	ated: C.O.C. No.: DUP
Purge Method: ☐ Centrifugal Pump ☑ Disposable Bailer ☒ Hand	ail □ Submersible Pump 🛛 Teflon Bailer □ Other
Purge Water Storage Container Type:	Storage Location: On Site
Date Purge Water Disposed:	Where Disposed: On Site
Analyses Requested No. and T	pe of Bottles Used 2.07 # 56.60
TPHg, TPHd, TPHmo	
BTEX, SVOC's, Lead Scavengers, Fuel Oxy's _	
Lab Name: Gurtis & Tompkins Test america	
Well NoM W - 7 Depth of Water	56.60
Well Diameter: Well Depth	
Ø 2" (0.16 gal/feet) □ 5" (1.02 gal/feet) Water Column He	10.35 80% DTW 58.67
□ 4" (0.65 gal/feet) □ 6" (1.47 gal/feet) Well Volume	80% DTW 38.6/
Inlet Denth Volume ORP 1	mnerature PH Cond DO

				T	1		•/3	1	T
Time	Inlet Depth	Depth to Water	Volume Purged (gal)	ORP	Temperature (C°)	PH (SU)	Cond (ne/cmc)	DO (mg/l)	Remarks
10:30 10:43 10:48			S	ta	V+				
10:43			1.75	41.3	64.87	6.59	.714	2.5	
10:48		3.5	3-50	18.4	6 4.83	\$.68	-719	2.10	
10:57		5.25	1. 25	26.0	64.78	6.67	.717	7.08	
11:10		7 -7	78.0	20.9 mp10	6598	6.76	.713	2.42	
11:15		56.70	Şa	nP10					
			/						
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								/	
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Project No. 001-09567-07-***	Date: March <u>17</u> , 2009	Page 1 of
Project Name <u>: Hanson Radum</u>	Sampling Location: Hans	of Radum
Sampler's Name: Tom Colling	Sample	No.: MW-8 DFB
Sampling Plan By: Ron Golobow		<u>^</u>
Purge Method: Centrifugal Pump Disposa	ıble Bailer 💆 Hand Bail 🏻 Submersible Pump 🕰 Tefl	lon Bailer □ Other
Purge Water Storage Container Type:		
Date Purge Water Disposed:	Where Disposed: On Site	<u> </u>
Analyses Requested	No. and Type of Bottles Used	1.926 + 54.67
TPHg, TPHd, TPHmo	•	
BTEX, SVÓC's, Lead Scavengers, Fuel Oxy's _		
Lab Name: Gurtis & Tompkins 1651	America	
Delivery By	X Hand	
Well No. MW-8	Depth of Water 54.67	
Well Diameter:	Well Depth 44.30	
★ 2" (0.16 gal/feet) □ 5" (1.02 gal/feet)	Water Çolumn Height	
☐ 4" (0.65 gal/feet) ☐ 6" (1.47 gal/feet)	Well Volume 1.541 ~ 1.5	80% DTW <u>56.596</u>

Time	Inlet Depth	Depth to Water	Volume Purged (gal)	ORP	Temperature (C°)	PH (SU)	Cond (us/cmc)	DO (mg/l)	Remarks	
14:50				Staf	- Purz	e			well be	185
12:05			1.5	33.6	67.24	6.43	.923		LOCK	cal De
12:12			3.0	-2.5	66.07	7.07	.931	1.94		
12:18			4.5	-2.8	66.07	7.21	.921	1.59		40.74
12:23			6.0	4.7	66.38	7.20	.921	2.25		
12:30	54.	75 S		11e						
12:45		D	UP							
1										
				-Tm						
			-	T.K						
								/		

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Project No. 001-09567-07-***	Date: March <u>1</u>	, 2009	Page 1 of
Project Name <u>: Hanson Radum</u>	Sampling Location	n: Haffon Rady	
Sampler's Name: A. Gill KOCTI			
Sampling Plan By: Ron Golobow		C.O.C. No.:	
Purge Method: ☐ Centrifugal Pump ☐ Disposable Baile	er □ Hand Bail □ Submersible	Pump ☐ Teflon Bailer ☐ Other	
Purge Water Storage Container Type:	Storage Location	on: On Site	
Date Purge Water Disposed:	Where Dispose	ed: On Site	
Analyses Requested	No. and Type of Bottles Used	····	
TPHg, TPHd, TPHmo			2) = .912
BTEX, SVOC's, Lead Scavengers, Fuel Oxy's	<u> </u>	——————————————————————————————————————	
Lab Name: Curtis & Tompkins TEST ame	ricu		
Delivery By ☐ Courier ➡ Har	nd		
Well No. Put Q Depth	of Water <u>55.51</u>		
- 18	Depth <u>55.07</u>		
,	Column Height 4,56		0.00
• • • • • • • • • • • • • • • • • • • •	/olume <u>* 73 へ.75</u>	(€ WTG %08 I	1.7 44

Time	Inlet Depth	Depth to Water	Volume Purged (gal)	ORP	Temperature (C°)	PH (SU)	. Cond (μs/cmc)	DO (mg/l)	Remarks
R:3b			•75	,9	7.08	7.05	795	1.62	and the second s
845			1.5	-1.6	:7.74	7.08	796	1.93	
950 950			a-a5	-39	17.95	7.08	803	ରୁ. ୮ର	
853			3	-7.9	18.05	7.01	පිටපි	a.44	
258			3.75	-9.9	17.82	7.05	806	a.13	
902			4.5	-11.6	18.12	7.05	810	9.91	
906			5.35	-21.9	17.42	7.09	810	३.७०	
911			6	-33.4	18.19	7.06	821	a·57	
13916			6.75	-43.2	17.69	7.08	813	2.30	
0920			7.5	-67.9	18.18	7.07	822	a. 39	
0926			8,25	-78.6	17.91	7.05	815	à.08	
0930			9	-79.7	17.97	7,07	813	2.53	·
0942		50.52	975	Scur	, de				
					(RC))	-		·

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WATER-QUALITY SAMPLING LOG

Project No. 001-09567-07-***	Date: March 17, 2009		age 1 of
Project Name <u>: Hanson Radum</u>	Sampling Location: Ha	5506 Radun	
Project Name: Hanson Radum Sampler's Name: A Gilveatte	Sam	ple No.: MW-10	□ FB
Sampling Plan By: Ron Golobow	Dated:	C.O.C. No.:	□ DUP
Purge Method: Centrifugal Pump Disposabl	e Bailer 🗷 Hand Bail 🛘 Submersible Pump 🗹	Teflon Bailer ☐ Other	
Purge Water Storage Container Type:	Storage Location: On Si	te	
Date Purge Water Disposed:	Where Disposed: On	Site	
Analyses Requested	No. and Type of Bottles Used		
TPHg, TPHd, TPHmo		6	
BTEX, SVOC's, Lead Scavengers, Fuel Oxy's _		502	11
Lab Name: Gurtis & Tempkins Test ameri	4	1,32+ 50.3	ч
Delivery By	Hand		
Well Diameter: 5" (1.02 gal/feet)	Nell Depth 56.93 Water Column Height 6.59 Water Volume 1.05	80% DTW <u>51.6</u> 6	0

Time	Inlet Depth	Depth to Water	Volume Purged (gal)	ORP	Temperature (C°)	PH (SU)	Cond (µs/cmc)	DO (mg/l)	Remarks
3 1425									Start purging
1430	S			-2.21.7	19.18	6.97	1099	2.68	
1432				-221.6	19.00	6.88	1086	2.68	
1436				-217.5	18.58	6.87	1075	1.83	
1441				-184.8	19.12	6.87	1084	a.72	
1449		50.54							Sample
	_							·	,
				/					
			/	11	(U)				
	.,								

Continue remarks on reverse, if needed.

APPENDIX B

Laboratory Certified Analytical Reports



ANALYTICAL REPORT

Job Number: 720-18564-1

Job Description: Hanson Radum

For: LFR, Inc. 1900 Powell St 12th Floor Emeryville, CA 94608-1827

Attention: Mr. Ron Goloubow

Approved for release Afsaneh Salimpour Project Manager I 3/30/2009 5:16 PM

Afsaneh Salimpour Project Manager I afsaneh.salimpour@testamericainc.com 03/30/2009

Akanef Sal

Job Narrative 720-J18564-1

Comments

No additional comments.

Receipt

Received 3 voas labeled MW-9. 1 of the 3 voas the time on it is 12:45. Time on the other 2 voas match the COC:9:42. The sample that is labeled for 12:45 looks very different than the other 2. Labeled as MW-9 and wrote on container for the lab not to use this voa.

All other samples were received in good condition within temperature requirements.

GC/MS VOA

Method(s) 8260B/CA_LUFTMS: The matrix spike / matrix spike duplicate (MS/MSD) recoveries for batch 47947 were outside control limits. The associated laboratory control sample (LCS) met acceptance criteria.

No other analytical or quality issues were noted.

GC/MS Semi VOA

No analytical or quality issues were noted.

GC Semi VOA

No analytical or quality issues were noted.

Organic Prep

Method(s) 3510C SGC: not enough sample; only one amber provided for diesel .

No other analytical or quality issues were noted.

EXECUTIVE SUMMARY - Detections

Client: LFR, Inc. Job Number: 720-18564-1

Lab Sample ID Client Sample ID Reporting
Analyte Result / Qualifier Limit Units Method

No Detections

METHOD SUMMARY

Client: LFR, Inc. Job Number: 720-18564-1

Description	Lab Location	Method	Preparation Method
Matrix: Water			
Volatile Organic Compounds by GC/MS Purge and Trap	TAL SF TAL SF	SW846 8260B	/CA_LUFTMS SW846 5030B
Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS) Liquid-Liquid Extraction (Separatory Funnel)	TAL SF	SW846 8270C	SW846 3510C
Diesel Range Organics (DRO) (GC) Liquid-Liquid Extraction (Separatory Funnel) Liquid-Liquid Extraction (Separatory Funnel)	TAL SF TAL SF TAL SF	SW846 8015B	SW846 3510C SW846 3510C SGC

Lab References:

TAL SF = TestAmerica San Francisco

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

METHOD / ANALYST SUMMARY

Client: LFR, Inc. Job Number: 720-18564-1

Method	Analyst	Analyst ID
SW846 8260B/CA_LUFTMS	Ali, Badri	ВА
SW846 8270C	Lee, Michael	ML
SW846 8015B SW846 8015B	Hayashi, Derek Relja, Marlene	DH MR

SAMPLE SUMMARY

Client: LFR, Inc. Job Number: 720-18564-1

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received
720-18564-1	MW-5	Water	03/17/2009 0935	03/17/2009 1545
720-18564-2	MW-9	Water	03/17/2009 0942	03/17/2009 1545
720-18564-3	MW-7	Water	03/17/2009 1115	03/17/2009 1545
720-18564-4	MW-6	Water	03/17/2009 1055	03/17/2009 1545
720-18564-5	MW-8	Water	03/17/2009 1230	03/17/2009 1545
720-18564-6	MW-2	Water	03/17/2009 1212	03/17/2009 1545
720-18564-7	MW-10	Water	03/17/2009 1449	03/17/2009 1545
720-18564-8FD	MW-8-DUP	Water	03/17/2009 1245	03/17/2009 1545
720-18564-9FB	FB-031709	Water	03/17/2009 0830	03/17/2009 1545

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-5

 Lab Sample ID:
 720-18564-1
 Date Sampled:
 03/17/2009 0935

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA_LUFTMS Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Initial Weight/Volume: 10 mL Date Analyzed: 03/23/2009 1454 Final Weight/Volume: 10 mL

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	90		78 - 112
1,2-Dichloroethane-d4 (Surr)	87		67 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-9

 Lab Sample ID:
 720-18564-2
 Date Sampled:
 03/17/2009 0942

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA_LUFTMS Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Initial Weight/Volume: 10 mL Date Analyzed: 03/23/2009 1648 Final Weight/Volume: 10 mL

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	92		78 - 112
1,2-Dichloroethane-d4 (Surr)	92		67 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-7

 Lab Sample ID:
 720-18564-3
 Date Sampled:
 03/17/2009 1115

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA_LUFTMS Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Initial Weight/Volume: 10 mL Date Analyzed: 03/23/2009 1711 Final Weight/Volume: 10 mL

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	89		78 - 112
1,2-Dichloroethane-d4 (Surr)	92		67 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-6

Lab Sample ID: 720-18564-4 Date Sampled: 03/17/2009 1055 Client Matrix: Water Date Received: 03/17/2009 1545

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA_LUFTMS Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Initial Weight/Volume: 10 mL 10 mL Date Analyzed: 03/23/2009 1733 Final Weight/Volume:

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	90		78 - 112
1.2-Dichloroethane-d4 (Surr)	93		67 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-8

 Lab Sample ID:
 720-18564-5
 Date Sampled:
 03/17/2009
 1230

 Client Matrix:
 Water
 Date Received:
 03/17/2009
 1545

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA_LUFTMS Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Initial Weight/Volume: 10 mL Date Analyzed: 03/23/2009 1756 Final Weight/Volume: 10 mL

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec	Acceptance	e Limits
Toluene-d8 (Surr)	92	78 - 112	
1,2-Dichloroethane-d4 (Surr)	95	67 - 126	

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-2

Lab Sample ID: 720-18564-6 Date Sampled: 03/17/2009 1212 Client Matrix: Water Date Received: 03/17/2009 1545

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA_LUFTMS Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Initial Weight/Volume: 10 mL 10 mL Date Analyzed: 03/23/2009 1819 Final Weight/Volume:

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	92		78 - 112
1.2-Dichloroethane-d4 (Surr)	90		67 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-10

 Lab Sample ID:
 720-18564-7
 Date Sampled:
 03/17/2009 1449

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA_LUFTMS Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Initial Weight/Volume: 10 mL Date Analyzed: 03/23/2009 1841 Final Weight/Volume: 10 mL

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	91		78 - 112
1,2-Dichloroethane-d4 (Surr)	91		67 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-8-DUP

 Lab Sample ID:
 720-18564-8FD
 Date Sampled:
 03/17/2009 1245

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA LUFTMS Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Initial Weight/Volume: 10 mL Date Analyzed: 03/23/2009 1904 Final Weight/Volume: 10 mL

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	96		78 - 112
1,2-Dichloroethane-d4 (Surr)	94		67 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: FB-031709

 Lab Sample ID:
 720-18564-9FB
 Date Sampled:
 03/17/2009 0830

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA_LUFTMS Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Initial Weight/Volume: 10 mL Date Analyzed: 03/23/2009 1431 Final Weight/Volume: 10 mL

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	90		78 - 112
1,2-Dichloroethane-d4 (Surr)	87		67 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-5

 Lab Sample ID:
 720-18564-1
 Date Sampled:
 03/17/2009 0935

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

Method: 8270C Analysis Batch: 720-48005 Instrument ID: Sat 2K

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-

Dilution: 1.0 Initial Weight/Volume: 980 mL

Date Analyzed: 03/25/2009 1403 Final Weight/Volume: 1 mL

Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Phenol	ND		2.0
Bis(2-chloroethyl)ether	ND		2.0
2-Chlorophenol	ND		2.0
1,3-Dichlorobenzene	ND		2.0
1,4-Dichlorobenzene	ND		2.0
Benzyl alcohol	ND		5.1
1,2-Dichlorobenzene	ND		2.0
2-Methylphenol	ND		2.0
4-Methylphenol	ND		2.0
N-Nitrosodi-n-propylamine	ND		2.0
Hexachloroethane	ND		2.0
Nitrobenzene	ND		2.0
Isophorone	ND		2.0
2-Nitrophenol	ND		2.0
2,4-Dimethylphenol	ND		2.0
Bis(2-chloroethoxy)methane	ND		5.1
2,4-Dichlorophenol	ND		5.1
1,2,4-Trichlorobenzene	ND		2.0
Naphthalene	ND		2.0
4-Chloroaniline	ND		2.0
Hexachlorobutadiene	ND		2.0
4-Chloro-3-methylphenol	ND		5.1
2-Methylnaphthalene	ND		2.0
Hexachlorocyclopentadiene	ND		5.1
2,4,6-Trichlorophenol	ND		2.0
2,4,5-Trichlorophenol	ND		2.0
2-Chloronaphthalene	ND		2.0
2-Nitroaniline	ND		10
Dimethyl phthalate	ND		5.1
Acenaphthylene	ND		2.0
3-Nitroaniline	ND		5.1
Acenaphthene	ND		2.0
2,4-Dinitrophenol	ND		10
4-Nitrophenol	ND		10
Dibenzofuran	ND		2.0
2,4-Dinitrotoluene	ND		2.0
2,6-Dinitrotoluene	ND		5.1
Diethyl phthalate	ND		5.1
4-Chlorophenyl phenyl ether	ND		5.1
Fluorene	ND		2.0
4-Nitroaniline	ND		10
2-Methyl-4,6-dinitrophenol	ND		10
N-Nitrosodiphenylamine	ND		2.0
4-Bromophenyl phenyl ether	ND		5.1

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-5

Lab Sample ID: 720-18564-1 Date Sampled: 03/17/2009 0935 Client Matrix: Water 03/17/2009 1545 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Dilution: 1.0

Initial Weight/Volume: 980 mL Date Analyzed: 03/25/2009 1403 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.0
Pentachlorophenol	ND		10
Phenanthrene	ND		2.0
Anthracene	ND		2.0
Di-n-butyl phthalate	ND		5.1
Fluoranthene	ND		2.0
Pyrene	ND		2.0
Butyl benzyl phthalate	ND		5.1
3,3'-Dichlorobenzidine	ND		5.1
Benzo[a]anthracene	ND		5.1
Bis(2-ethylhexyl) phthalate	ND		10
Chrysene	ND		2.0
Di-n-octyl phthalate	ND		20
Benzo[b]fluoranthene	ND		2.0
Benzo[a]pyrene	ND		2.0
Benzo[k]fluoranthene	ND		2.0
Indeno[1,2,3-cd]pyrene	ND		2.0
Benzo[g,h,i]perylene	ND		2.0
Benzoic acid	ND		10
Azobenzene	ND		2.0
Dibenz(a,h)anthracene	ND		2.0
Surrogate	%Rec		Acceptance Limits
Nitrobenzene-d5	69		6 - 98
2-Fluorobiphenyl	68		6 - 103
Terphenyl-d14	75		36 - 106
2-Fluorophenol	40		1 - 66
Phenol-d5	27		1 - 47
2,4,6-Tribromophenol	70		22 - 124

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-9

Lab Sample ID: 720-18564-2 Date Sampled: 03/17/2009 0942 Client Matrix: Water 03/17/2009 1545 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-Initial Weight/Volume: 950 mL

Dilution: 1.0

Date Analyzed: 03/25/2009 1436 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Phenol	ND		2.1
Bis(2-chloroethyl)ether	ND		2.1
2-Chlorophenol	ND		2.1
1,3-Dichlorobenzene	ND		2.1
1,4-Dichlorobenzene	ND		2.1
Benzyl alcohol	ND		5.3
1,2-Dichlorobenzene	ND		2.1
2-Methylphenol	ND		2.1
4-Methylphenol	ND		2.1
N-Nitrosodi-n-propylamine	ND		2.1
Hexachloroethane	ND		2.1
Nitrobenzene	ND		2.1
Isophorone	ND		2.1
2-Nitrophenol	ND		2.1
2,4-Dimethylphenol	ND		2.1
Bis(2-chloroethoxy)methane	ND		5.3
2,4-Dichlorophenol	ND		5.3
1,2,4-Trichlorobenzene	ND		2.1
Naphthalene	ND		2.1
4-Chloroaniline	ND		2.1
Hexachlorobutadiene	ND		2.1
4-Chloro-3-methylphenol	ND		5.3
2-Methylnaphthalene	ND		2.1
Hexachlorocyclopentadiene	ND		5.3
2,4,6-Trichlorophenol	ND		2.1
2,4,5-Trichlorophenol	ND		2.1
2-Chloronaphthalene	ND		2.1
2-Nitroaniline	ND		11
Dimethyl phthalate	ND		5.3
Acenaphthylene	ND		2.1
3-Nitroaniline	ND		5.3
Acenaphthene	ND		2.1
2,4-Dinitrophenol	ND		11
4-Nitrophenol	ND		11
Dibenzofuran	ND		2.1
2,4-Dinitrotoluene	ND		2.1
2,6-Dinitrotoluene	ND		5.3
Diethyl phthalate	ND		5.3
4-Chlorophenyl phenyl ether	ND		5.3
Fluorene	ND		2.1
4-Nitroaniline	ND		11
2-Methyl-4,6-dinitrophenol	ND		11
N-Nitrosodiphenylamine	ND		2.1
4-Bromophenyl phenyl ether	ND		5.3

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-9

Lab Sample ID: 720-18564-2 Date Sampled: 03/17/2009 0942 Client Matrix: Water 03/17/2009 1545 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Dilution: 1.0

Initial Weight/Volume: 950 mL Date Analyzed: 03/25/2009 1436 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.1
Pentachlorophenol	ND		11
Phenanthrene	ND		2.1
Anthracene	ND		2.1
Di-n-butyl phthalate	ND		5.3
Fluoranthene	ND		2.1
Pyrene	ND		2.1
Butyl benzyl phthalate	ND		5.3
3,3'-Dichlorobenzidine	ND		5.3
Benzo[a]anthracene	ND		5.3
Bis(2-ethylhexyl) phthalate	ND		11
Chrysene	ND		2.1
Di-n-octyl phthalate	ND		21
Benzo[b]fluoranthene	ND		2.1
Benzo[a]pyrene	ND		2.1
Benzo[k]fluoranthene	ND		2.1
Indeno[1,2,3-cd]pyrene	ND		2.1
Benzo[g,h,i]perylene	ND		2.1
Benzoic acid	ND		11
Azobenzene	ND		2.1
Dibenz(a,h)anthracene	ND		2.1
Surrogate	%Rec		Acceptance Limits
Nitrobenzene-d5	67		6 - 98
2-Fluorobiphenyl	63		6 - 103
Terphenyl-d14	73		36 - 106
2-Fluorophenol	40		1 - 66
Phenol-d5	27		1 - 47
2,4,6-Tribromophenol	77		22 - 124

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-7

 Lab Sample ID:
 720-18564-3
 Date Sampled:
 03/17/2009 1115

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

Method: 8270C Analysis Batch: 720-48005 Instrument ID: Sat 2K²

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-

 Dilution:
 1.0
 Initial Weight/Volume:
 980 mL

 Date Analyzed:
 03/25/2009 1510
 Final Weight/Volume:
 1 mL

 Date Prepared:
 03/23/2009 1257
 Injection Volume:
 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Phenol	ND		2.0
Bis(2-chloroethyl)ether	ND		2.0
2-Chlorophenol	ND		2.0
1,3-Dichlorobenzene	ND		2.0
1,4-Dichlorobenzene	ND		2.0
Benzyl alcohol	ND		5.1
1,2-Dichlorobenzene	ND		2.0
2-Methylphenol	ND		2.0
4-Methylphenol	ND		2.0
N-Nitrosodi-n-propylamine	ND		2.0
Hexachloroethane	ND		2.0
Nitrobenzene	ND		2.0
Isophorone	ND		2.0
2-Nitrophenol	ND		2.0
2,4-Dimethylphenol	ND		2.0
Bis(2-chloroethoxy)methane	ND		5.1
2,4-Dichlorophenol	ND		5.1
1,2,4-Trichlorobenzene	ND		2.0
Naphthalene	ND		2.0
4-Chloroaniline	ND		2.0
Hexachlorobutadiene	ND		2.0
4-Chloro-3-methylphenol	ND		5.1
2-Methylnaphthalene	ND		2.0
Hexachlorocyclopentadiene	ND		5.1
2,4,6-Trichlorophenol	ND		2.0
2,4,5-Trichlorophenol	ND		2.0
2-Chloronaphthalene	ND		2.0
2-Nitroaniline	ND		10
Dimethyl phthalate	ND		5.1
Acenaphthylene	ND		2.0
3-Nitroaniline	ND		5.1
Acenaphthene	ND		2.0
2,4-Dinitrophenol	ND		10
4-Nitrophenol	ND		10
Dibenzofuran	ND		2.0
2,4-Dinitrotoluene	ND		2.0
2,6-Dinitrotoluene	ND		5.1
Diethyl phthalate	ND		5.1
4-Chlorophenyl phenyl ether	ND		5.1
Fluorene	ND		2.0
4-Nitroaniline	ND		10
2-Methyl-4,6-dinitrophenol	ND		10
N-Nitrosodiphenylamine	ND		2.0
4-Bromophenyl phenyl ether	ND		5.1

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-7

Lab Sample ID: 720-18564-3 Date Sampled: 03/17/2009 1115 Client Matrix: Water Date Received: 03/17/2009 1545

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-

Dilution: 1.0

Initial Weight/Volume: 980 mL Final Weight/Volume: Date Analyzed: 03/25/2009 1510 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.0
Pentachlorophenol	ND		10
Phenanthrene	ND		2.0
Anthracene	ND		2.0
Di-n-butyl phthalate	ND		5.1
Fluoranthene	ND		2.0
Pyrene	ND		2.0
Butyl benzyl phthalate	ND		5.1
3,3'-Dichlorobenzidine	ND		5.1
Benzo[a]anthracene	ND		5.1
Bis(2-ethylhexyl) phthalate	ND		10
Chrysene	ND		2.0
Di-n-octyl phthalate	ND		20
Benzo[b]fluoranthene	ND		2.0
Benzo[a]pyrene	ND		2.0
Benzo[k]fluoranthene	ND		2.0
Indeno[1,2,3-cd]pyrene	ND		2.0
Benzo[g,h,i]perylene	ND		2.0
Benzoic acid	ND		10
Azobenzene	ND		2.0
Dibenz(a,h)anthracene	ND		2.0
Surrogate	%Rec		Acceptance Limits
Nitrobenzene-d5	76		6 - 98
2-Fluorobiphenyl	75		6 - 103
Terphenyl-d14	85		36 - 106
2-Fluorophenol	45		1 - 66
Phenol-d5	30		1 - 47
2,4,6-Tribromophenol	82		22 - 124

1.0 uL

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-6

03/23/2009 1257

Date Prepared:

 Lab Sample ID:
 720-18564-4
 Date Sampled:
 03/17/2009 1055

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

Method: 8270C Analysis Batch: 720-48005 Instrument ID: Sat 2K

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-

Injection Volume:

Dilution: 1.0 Initial Weight/Volume: 950 mL Date Analyzed: 03/25/2009 1544 Final Weight/Volume: 1 mL

Result (ug/L) Qualifier RL Analyte Phenol ND 2.1 ND Bis(2-chloroethyl)ether 2.1 2-Chlorophenol ND 2.1 1,3-Dichlorobenzene 2.1 ND 1,4-Dichlorobenzene ND 2.1 Benzvl alcohol ND 5.3 1,2-Dichlorobenzene ND 2.1 2-Methylphenol ND 2.1 4-Methylphenol ND 2.1 N-Nitrosodi-n-propylamine ND 2.1 Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Isophorone ND 2.1 2-Nitrophenol 2.1 ND 2,4-Dimethylphenol ND 2.1 Bis(2-chloroethoxy)methane ND 5.3 2,4-Dichlorophenol ND 5.3 1,2,4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 Hexachlorobutadiene ND 2.1 4-Chloro-3-methylphenol ND 5.3 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 5.3 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.3 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.3 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.3 Diethyl phthalate ND 5.3 4-Chlorophenyl phenyl ether ND 5.3 Fluorene ND 2.1 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.1 4-Bromophenyl phenyl ether ND 5.3

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-6

Lab Sample ID: 720-18564-4 Date Sampled: 03/17/2009 1055 Client Matrix: Water 03/17/2009 1545 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Dilution: 1.0

Initial Weight/Volume: 950 mL Date Analyzed: 03/25/2009 1544 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.1
Pentachlorophenol	ND		11
Phenanthrene	ND		2.1
Anthracene	ND		2.1
Di-n-butyl phthalate	ND		5.3
Fluoranthene	ND		2.1
Pyrene	ND		2.1
Butyl benzyl phthalate	ND		5.3
3,3'-Dichlorobenzidine	ND		5.3
Benzo[a]anthracene	ND		5.3
Bis(2-ethylhexyl) phthalate	ND		11
Chrysene	ND		2.1
Di-n-octyl phthalate	ND		21
Benzo[b]fluoranthene	ND		2.1
Benzo[a]pyrene	ND		2.1
Benzo[k]fluoranthene	ND		2.1
Indeno[1,2,3-cd]pyrene	ND		2.1
Benzo[g,h,i]perylene	ND		2.1
Benzoic acid	ND		11
Azobenzene	ND		2.1
Dibenz(a,h)anthracene	ND		2.1
Surrogate	%Rec	A	cceptance Limits
Nitrobenzene-d5	53		6 - 98
2-Fluorobiphenyl	51	(6 - 103
Terphenyl-d14	66	;	36 - 106
2-Fluorophenol	30		1 - 66
Phenol-d5	20	•	1 - 47
2,4,6-Tribromophenol	52	2	22 - 124

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-8

 Lab Sample ID:
 720-18564-5
 Date Sampled:
 03/17/2009 1230

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

Method: 8270C Analysis Batch: 720-48005 Instrument ID: Sat 2K

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-

Dilution: 1.0 Initial Weight/Volume: 980 mL
Date Analyzed: 03/25/2009 1617 Final Weight/Volume: 1 mL

 Date Analyzed:
 03/25/2009 1617
 Final Weight/Volume:
 1 mL

 Date Prepared:
 03/23/2009 1257
 Injection Volume:
 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Phenol	ND		2.0
Bis(2-chloroethyl)ether	ND		2.0
2-Chlorophenol	ND		2.0
1,3-Dichlorobenzene	ND		2.0
1,4-Dichlorobenzene	ND		2.0
Benzyl alcohol	ND		5.1
1,2-Dichlorobenzene	ND		2.0
2-Methylphenol	ND		2.0
4-Methylphenol	ND		2.0
N-Nitrosodi-n-propylamine	ND		2.0
Hexachloroethane	ND		2.0
Nitrobenzene	ND		2.0
Isophorone	ND		2.0
2-Nitrophenol	ND		2.0
2,4-Dimethylphenol	ND		2.0
Bis(2-chloroethoxy)methane	ND		5.1
2,4-Dichlorophenol	ND		5.1
1,2,4-Trichlorobenzene	ND		2.0
Naphthalene	ND		2.0
4-Chloroaniline	ND		2.0
Hexachlorobutadiene	ND		2.0
4-Chloro-3-methylphenol	ND		5.1
2-Methylnaphthalene	ND		2.0
Hexachlorocyclopentadiene	ND		5.1
2,4,6-Trichlorophenol	ND		2.0
2,4,5-Trichlorophenol	ND		2.0
2-Chloronaphthalene	ND		2.0
2-Nitroaniline	ND		10
Dimethyl phthalate	ND		5.1
Acenaphthylene	ND		2.0
3-Nitroaniline	ND		5.1
Acenaphthene	ND		2.0
2,4-Dinitrophenol	ND		10
4-Nitrophenol	ND		10
Dibenzofuran	ND		2.0
2,4-Dinitrotoluene	ND		2.0
2,6-Dinitrotoluene	ND		5.1
Diethyl phthalate	ND		5.1
4-Chlorophenyl phenyl ether	ND		5.1
Fluorene	ND		2.0
4-Nitroaniline	ND		10
2-Methyl-4,6-dinitrophenol	ND		10
N-Nitrosodiphenylamine	ND		2.0
4-Bromophenyl phenyl ether	ND		5.1

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: **MW-8**

Lab Sample ID: 720-18564-5 Date Sampled: 03/17/2009 1230 Client Matrix: 03/17/2009 1545 Water Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Dilution: 1.0

Initial Weight/Volume: 980 mL Date Analyzed: 03/25/2009 1617 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.0
Pentachlorophenol	ND		10
Phenanthrene	ND		2.0
Anthracene	ND		2.0
Di-n-butyl phthalate	ND		5.1
Fluoranthene	ND		2.0
Pyrene	ND		2.0
Butyl benzyl phthalate	ND		5.1
3,3'-Dichlorobenzidine	ND		5.1
Benzo[a]anthracene	ND		5.1
Bis(2-ethylhexyl) phthalate	ND		10
Chrysene	ND		2.0
Di-n-octyl phthalate	ND		20
Benzo[b]fluoranthene	ND		2.0
Benzo[a]pyrene	ND		2.0
Benzo[k]fluoranthene	ND		2.0
Indeno[1,2,3-cd]pyrene	ND		2.0
Benzo[g,h,i]perylene	ND		2.0
Benzoic acid	ND		10
Azobenzene	ND		2.0
Dibenz(a,h)anthracene	ND		2.0
Surrogate	%Rec		Acceptance Limits
Nitrobenzene-d5	63		6 - 98
2-Fluorobiphenyl	63		6 - 103
Terphenyl-d14	71		36 - 106
2-Fluorophenol	35		1 - 66
Phenol-d5	24		1 - 47
2,4,6-Tribromophenol	71		22 - 124

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-2

 Lab Sample ID:
 720-18564-6
 Date Sampled:
 03/17/2009 1212

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

Method: 8270C Analysis Batch: 720-48005 Instrument ID: Sat 2K²

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-

Dilution: 1.0 Initial Weight/Volume: 910 mL

Date Analyzed: 03/25/2009 1650 Final Weight/Volume: 1 mL

Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Bis(2-chlorophenol ND 2.2	Analyte	Result (ug/L)	Qualifier	RL
2-Chlorophenof ND 2.2 1,4-Dichlorobenzene ND 2.2 1,4-Dichlorobenzene ND 2.2 Benzyl alcohol ND 2.2 Benzyl alcohol ND 2.2	Phenol			2.2
1,3-Dichlorobenzene ND 2.2 Benzyl alcohol ND 5.5 1,2-Dichlorobenzene ND 5.5 2.2 Methylphenol ND 2.2 4-Methylphenol ND 2.2 4-Methylphenol ND 2.2 4-Methylphenol ND 2.2 Hexachloroethane ND 2.2 Hitrobenzene ND 2.2 Illopheroene ND 2.2 Nitrobenzene ND 2.2 2-Nitrophenol ND 2.2 2-Nitrophenol ND 2.2 2-Nitrophenol ND 2.2 2-4-Dienthylphenol ND 5.5 2-4-Dichlorophenol ND 5.5 2-4-Dichlorophenol ND 5.5 2-4-Dichlorophenol ND 2.2 Naphthalene ND 2.2 4-Chloro-3-methylphenol ND 2.2 Hexachlorobutadiene ND 2.2 Hexachlorophenol ND 2.2 Hexachlorophenol ND 2.2 H	Bis(2-chloroethyl)ether	ND		
1.4-Dichlorobenzene ND 5.5 1.2-Dichlorobenzene ND 2.2 2-Methylphenol ND 2.2 4-Methylphenol ND 2.2 N-Nitrosodi-n-propylamine ND 2.2 N-Nitrosedi-n-propylamine ND 2.2 N-Nitrobenzene ND 2.2 Isophorone ND 2.2 2-Nitrophenol ND 5.5 1,2-4-Tichlorobenzene ND 5.5 ND 2.2 4-Chlorosalidiene ND 2.2 4-Chlorosalidiene ND 2.2 4-Chlorosalidiene ND 2.2 4-Kaschlorobutadiene ND 5.5 4-Kaschlorobutadiene ND 5.5 4-Kaschlorobutadiene ND 5.5 4-Kaschlorobutadiene ND 5.5 4-Kas	2-Chlorophenol	ND		
Benzyl alcohol ND 5.5 1,2-Dichlorobenzene ND 2.2 2-Methylphenol ND 2.2 4-Methylphenol ND 2.2 4-Neitrosodi-n-propylamine ND 2.2 Hexachloroethane ND 2.2 Nitrobenzene ND 2.2 Isophorone ND 2.2 2-Nitrophenol ND 2.2 2-Nitrophenol ND 2.2 2-Nitrophenol ND 2.2 2-Nitrophenol ND 5.5 2,4-Dichlorophenol ND 5.5 2,4-Dichlorophenol ND 2.2 Naphthalene ND 2.2 4-Chloro-3-methylphenol ND 2.2 4-Chloro-3-methylphenol ND 2.2 4-Chloro-3-methylphenol ND 5.5 2,4-5-Trichlorophenol ND 5.5 2,4-5-Trichlorophenol ND 2.2 2,4-5-Trichlorophenol ND 2.2 2-Nitroanline <t< td=""><td>1,3-Dichlorobenzene</td><td>ND</td><td></td><td></td></t<>	1,3-Dichlorobenzene	ND		
1,2-Dichlorobenzene ND 2,2 2-Methylphenol ND 2,2 4-Methylphenol ND 2,2 N-Nitrosodin-propylamine ND 2,2 N-Nitrosodin-propylamine ND 2,2 NItrobenzene ND 2,2 Nitrobenzene ND 2,2 Shitrobenzene ND 2,2 Shitrobenzene ND 2,2 2-Nitrophenol ND 2,2 2-Liphichlorophenol ND 2,2 2-Liphichlorophenol ND 5,5 2-Liphichlorophenol ND 2,2 4-Chiloro-Arrichlorophenol ND 2,2 4-Chloro-Arrichlorophenol ND 2,2 4-Chloro-Arrichlorophenol ND 2,2 2-Methylnaphthalene ND 2,2 2-Metrylnaphthalene ND 2,2 2-Mitropaphenol ND 2,2 2-Mitropaphenol ND 2,2 2-Nitroaniline ND 2,2 Nitroaniline	1,4-Dichlorobenzene	ND		
2-Methylphenol ND 2.2 4-Methylphenol ND 2.2 N-Nitrosodi-n-propylamine ND 2.2 Hexachloroethane ND 2.2 Nitrobenzene ND 2.2 Isophorone ND 2.2 2-Nitrophenol ND 2.2 2-Nitrophenol ND 2.2 2-4-Dinethylphenol ND 2.2 Bis(2-chloroethoxy)methane ND 5.5 2,4-Dichlorophenol ND 5.5 2,4-Dichlorophenol ND 5.5 2,4-Dichlorophenol ND 5.5 2,4-Dichlorophenol ND 2.2 Naphthalene ND 2.2 4-Chloro-3-methylphenol ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 5.5 2-Methylnaphthalene ND 2.2 4-Klorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND	Benzyl alcohol			
4-Methylphenol ND 2.2 N-Nitrosodi-n-propylamine ND 2.2 Hexachloroethane ND 2.2 Nitrobenzene ND 2.2 Isophorone ND 2.2 2-Nitrophenol ND 2.2 2-Nitrophenol ND 2.2 2-Nitrophenol ND 2.2 18is(2-chloroethoxy)methane ND 5.5 2,4-Dichlorophenol ND 5.5 1,2,4-Tichlorophenol ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 2.2 Hexachlorocyclopentadiene ND 2.2 Hexachlorocyclopentadiene ND 5.5 2,4-B-Trichlorophenol ND 2.2 2,4-S-Trichlorophenol ND 2.2 2,4-S-Trichlorophenol ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND	•			
N-Nitrosodi-n-propylamine ND Lexachloroethane ND Lexachloroethoxyjmethane Lexachloroethoxyjmethane ND Lexachloroethane ND Lexachlo				
Hexachloroethane ND 2.2 Nitrobenzene ND 2.2 Isophorone ND 2.2 2-Nitrophenol ND 2.2 2-Homethylphenol ND 2.2 18(2-chloroethoxy)methane ND 5.5 2,4-Dichlorophenol ND 5.5 1,2,4-Trichlorobenzene ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 5.5 2-Methylnaphthalene ND 5.5 2-Methylnaphthalene ND 5.5 2-4,6-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-Nitroaniline ND 2.2 2-Nitroaniline ND 5.5 Acenaphthylene ND 5.5 Acenaphthylene	4-Methylphenol			
Nitrobenzene ND 2.2 Isophorone ND 2.2 2-Nitrophenol ND 2.2 2-Nitrophenol ND 2.2 3-Dimethylphenol ND 2.2 Bis(2-chloroethoxy)methane ND 5.5 2,4-Dichlorophenol ND 5.5 1,2,4-Trichlorobenzene ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 2.2 4-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 5.5 2,4-G-Trichlorophenol ND 2.2 2,4-5-Trichlorophenol ND 2.2 2-4,1-1richlorophenol ND 2.2 2-Nitroaniline ND 2.2 2-Nitroaniline ND 5.5 Dimethyl phthalate ND 5.5 Acenaphthene ND 2.2 2-4-Dinitrophenol	N-Nitrosodi-n-propylamine			
Isophorone ND 2.2 2-Nitrophenol ND 2.2 2,4-Dimethylphenol ND 2.2 Bis(2-chloroethoxy)methane ND 5.5 2,4-Dichlorophenol ND 5.5 1,2,4-Trichlorophenzene ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 2.2 4-Chloro-3-methylphenol ND 2.2 4-Mexachlorocyclopentadiene ND 2.2 4-Mexachlorocyclopentadiene ND 2.2 2-Methylnaphthalene ND 2.2 2,4,6-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2,4-5-Trichlorophenol ND 2.2 2-Nitroaniline ND 2.2 2-Nitroaniline ND 5.5 Acenaphthylene ND 5.5 Acenaphthylene ND 11	Hexachloroethane	ND		
2-Nitrophenol ND 2.2 2,4-Dinethylphenol ND 5.5 2,4-Dichlorophenol ND 5.5 2,4-Dichlorophenol ND 5.5 1,2,4-Trichlorobenzene ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 5.5 4-Chloro-3-methylphenol ND 2.2 4-Kachlorocyclopentadiene ND 2.2 4-Methylnaphthalene ND 2.2 4-A,5-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2,4-5-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-Nitroaniline ND 2.2 2-Nitroaniline ND 5.5 Acenaphthylaphthalate ND 5.5 Acenaphthylap	Nitrobenzene			
2,4-Dimethylphenol ND 5.5 Bis(2-chloroethoxy)methane ND 5.5 2,4-Dichlorophenol ND 5.5 1,2,4-Trichlorobenzene ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 2.2 4-Ka-Trichlorophenol ND 2.2 2,4,6-Trichlorophenol ND 2.2 2-Khitroaniline ND 2.2 2-Chloronaphthalene ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 5.5 Acenaphthene ND 5.5 Acenaphthene ND 11 Valitrophenol ND 11 4-Nitrophenol ND 11 4-Nitrophenol ND 2.2 2,4-Dinitrotoluene ND 5.5	Isophorone	ND		2.2
Bis(2-chloroethoxy)methane ND 5.5 2,4-Dichlorophenol ND 5.5 1,2,4-Trichlorobenzene ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 5.5 2-Methylnaphthalene ND 5.5 2-Methylnaphthalene ND 5.5 2-4,6-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-Nitroaniline ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinin	2-Nitrophenol	ND		
2,4-Dichlorophenol ND 5.5 1,2,4-Trichlorobenzene ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 2.2 Hexachlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2-Nitroaniline ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 2.2 3-Nitroaniline ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5<	2,4-Dimethylphenol	ND		2.2
1,2,4-Trichlorobenzene ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 5.5 2,4,6-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalatene ND 11 2-Nitroaniline ND 5.5 Acenaphthylene ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 4-Nitrophenol ND 11 4-Nitrotoluene ND 2.2 2,4-Dinitrotoluene ND 5.5 2,6-Dinitrotoluene ND 5.5 3-Chilorophenyl phenyl ether ND 5.5	Bis(2-chloroethoxy)methane	ND		5.5
Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 2.2 2-4,6-Trichlorophenol ND 2.2 2-4,5-Trichlorophenol ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalatene ND 5.5 Acenaphthylene ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 5.5 3-6-Dinitrotoluene ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND<	2,4-Dichlorophenol	ND		5.5
4-Chloroanilline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 5.5 2,4,6-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2-Chloronaphthalene ND 2.2 2-Chloronaphthalene ND 11 Dimethyl phthalate ND 11 Acenaphtylene ND 5.5 Acenaphthylene ND 5.5 Acenaphthene ND 5.5 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 4-Nitrophenol ND 11 1-Indictoduene ND 2.2 2,4-Dinitrotoluene ND 5.5 2,6-Dinitrotoluene ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 5.5 4-Nitroaniline ND 5.5 -ND 5.5 5.5 <td>1,2,4-Trichlorobenzene</td> <td>ND</td> <td></td> <td>2.2</td>	1,2,4-Trichlorobenzene	ND		2.2
Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 5.5 2,4,6-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2-Chloronaphthalene ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 11 Acenaphthylene ND 5.5 Acenaphthylene ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 5.5 G-6-Dinitrotoluene ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 5.5 4-Nitrosoniline ND 11 2-Methyl-4,6-dinitrophenol	Naphthalene	ND		2.2
4-Chloro-3-methylphenol ND 5.5 2-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 5.5 2,4,6-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2-Chloronaphthalene ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 5.5 Acenaphthene ND 11 4-Nitrophenol ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 <t< td=""><td>4-Chloroaniline</td><td>ND</td><td></td><td>2.2</td></t<>	4-Chloroaniline	ND		2.2
2-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 5.5 2,4,6-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2-Chloronaphthalene ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 5.5 Acenaphthene ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 5.5 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	Hexachlorobutadiene	ND		2.2
Hexachlorocyclopentadiene ND 5.5 2,4,6-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2-Chloronaphthalene ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 5.5 Acenaphthene ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 11 4-Nitrophenol ND 2.2 2,4-Dinitrotoluene ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine <	4-Chloro-3-methylphenol	ND		5.5
Hexachlorocyclopentadiene ND 5.5 2,4,6-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2-Chloronaphthalene ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 5.5 Acenaphthene ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	2-Methylnaphthalene	ND		2.2
2,4,5-Trichlorophenol ND 2.2 2-Chloronaphthalene ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 11 2,4-Dinitrotoluene ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 NEW Johnstein ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 5.5 4-Nitroaniline ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	Hexachlorocyclopentadiene	ND		5.5
2-Chloronaphthalene ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 5.5 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	2,4,6-Trichlorophenol	ND		2.2
2-Nitroaniline ND 11 Dimethyl phthalate ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	2,4,5-Trichlorophenol	ND		2.2
Dimethyl phthalate ND 5.5 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 5.5 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	2-Chloronaphthalene	ND		2.2
Acenaphthylene ND 2.2 3-Nitroaniline ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	2-Nitroaniline	ND		11
3-Nitroaniline ND 5.5 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 11	Dimethyl phthalate	ND		5.5
Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	Acenaphthylene	ND		2.2
2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	3-Nitroaniline	ND		5.5
4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	Acenaphthene	ND		2.2
Dibenzofuran ND 2.2 2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	2,4-Dinitrophenol	ND		11
2,4-Dinitrotoluene ND 2.2 2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	4-Nitrophenol	ND		11
2,6-Dinitrotoluene ND 5.5 Diethyl phthalate ND 5.5 4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	Dibenzofuran	ND		2.2
Diethyl phthalateND5.54-Chlorophenyl phenyl etherND5.5FluoreneND2.24-NitroanilineND112-Methyl-4,6-dinitrophenolND11N-NitrosodiphenylamineND2.2	2,4-Dinitrotoluene	ND		
4-Chlorophenyl phenyl etherND5.5FluoreneND2.24-NitroanilineND112-Methyl-4,6-dinitrophenolND11N-NitrosodiphenylamineND2.2	2,6-Dinitrotoluene	ND		5.5
4-Chlorophenyl phenyl ether ND 5.5 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	Diethyl phthalate	ND		
Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2	4-Chlorophenyl phenyl ether	ND		5.5
2-Methyl-4,6-dinitrophenolND11N-NitrosodiphenylamineND2.2	Fluorene	ND		2.2
N-Nitrosodiphenylamine ND 2.2	4-Nitroaniline	ND		11
N-Nitrosodiphenylamine ND 2.2	2-Methyl-4,6-dinitrophenol	ND		11
	N-Nitrosodiphenylamine			
	4-Bromophenyl phenyl ether	ND		5.5

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-2

Lab Sample ID: 720-18564-6 Date Sampled: 03/17/2009 1212 Client Matrix: Water 03/17/2009 1545 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Dilution: 1.0

Initial Weight/Volume: 910 mL Date Analyzed: 03/25/2009 1650 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.2
Pentachlorophenol	ND		11
Phenanthrene	ND		2.2
Anthracene	ND		2.2
Di-n-butyl phthalate	ND		5.5
Fluoranthene	ND		2.2
Pyrene	ND		2.2
Butyl benzyl phthalate	ND		5.5
3,3'-Dichlorobenzidine	ND		5.5
Benzo[a]anthracene	ND		5.5
Bis(2-ethylhexyl) phthalate	ND		11
Chrysene	ND		2.2
Di-n-octyl phthalate	ND		22
Benzo[b]fluoranthene	ND		2.2
Benzo[a]pyrene	ND		2.2
Benzo[k]fluoranthene	ND		2.2
Indeno[1,2,3-cd]pyrene	ND		2.2
Benzo[g,h,i]perylene	ND		2.2
Benzoic acid	ND		11
Azobenzene	ND		2.2
Dibenz(a,h)anthracene	ND		2.2
Surrogate	%Rec	Ac	ceptance Limits
Nitrobenzene-d5	62	6	- 98
2-Fluorobiphenyl	59	6	- 103
Terphenyl-d14	70	3	6 - 106
2-Fluorophenol	38	1	- 66
Phenol-d5	27	1	- 47
2,4,6-Tribromophenol	66	2	2 - 124

RL

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-10

Lab Sample ID: 720-18564-7 Date Sampled: 03/17/2009 1449 Client Matrix: Water Date Received: 03/17/2009 1545

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

Method: 8270C Analysis Batch: 720-48005 Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Qualifier

Dilution: Initial Weight/Volume: 1.0 960 mL Date Analyzed: 03/25/2009 1724 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Result (ug/L) Analyte Phenol ND 2.1 ND Bis(2-chloroethyl)ether 2.1 2-Chlorophenol ND 2.1 1,3-Dichlorobenzene 2.1 ND 1,4-Dichlorobenzene ND 2.1 Benzvl alcohol ND 5.2 1,2-Dichlorobenzene ND 2.1 2-Methylphenol ND 2.1 4-Methylphenol ND 2.1 N-Nitrosodi-n-propylamine ND 2.1 Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Isophorone ND 2.1 2-Nitrophenol 2.1 ND 2,4-Dimethylphenol ND 2.1 Bis(2-chloroethoxy)methane ND 5.2 2,4-Dichlorophenol ND 5.2 1,2,4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 Hexachlorobutadiene ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2.1 2,4-Dinitrotoluene ND 5.2 2,6-Dinitrotoluene ND Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol 10 ND N-Nitrosodiphenylamine ND 2.1 4-Bromophenyl phenyl ether ND 5.2

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-10

Lab Sample ID: 720-18564-7 Date Sampled: 03/17/2009 1449 Client Matrix: Water 03/17/2009 1545 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Dilution: 1.0

Initial Weight/Volume: 960 mL Date Analyzed: 03/25/2009 1724 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.1
Pentachlorophenol	ND		10
Phenanthrene	ND		2.1
Anthracene	ND		2.1
Di-n-butyl phthalate	ND		5.2
Fluoranthene	ND		2.1
Pyrene	ND		2.1
Butyl benzyl phthalate	ND		5.2
3,3'-Dichlorobenzidine	ND		5.2
Benzo[a]anthracene	ND		5.2
Bis(2-ethylhexyl) phthalate	ND		10
Chrysene	ND		2.1
Di-n-octyl phthalate	ND		21
Benzo[b]fluoranthene	ND		2.1
Benzo[a]pyrene	ND		2.1
Benzo[k]fluoranthene	ND		2.1
Indeno[1,2,3-cd]pyrene	ND		2.1
Benzo[g,h,i]perylene	ND		2.1
Benzoic acid	ND		10
Azobenzene	ND		2.1
Dibenz(a,h)anthracene	ND		2.1
Surrogate	%Rec		Acceptance Limits
Nitrobenzene-d5	74		6 - 98
2-Fluorobiphenyl	69		6 - 103
Terphenyl-d14	76		36 - 106
2-Fluorophenol	44		1 - 66
Phenol-d5	32		1 - 47
2,4,6-Tribromophenol	75		22 - 124

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-8-DUP

Lab Sample ID: 720-18564-8FD Date Sampled: 03/17/2009 1245 Client Matrix: Water Date Received: 03/17/2009 1545

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID: 970 mL

Dilution: 1.0

Initial Weight/Volume: Date Analyzed: 03/25/2009 1757 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Phenol	ND		2.1
Bis(2-chloroethyl)ether	ND		2.1
2-Chlorophenol	ND		2.1
1,3-Dichlorobenzene	ND		2.1
1,4-Dichlorobenzene	ND		2.1
Benzyl alcohol	ND		5.2
1,2-Dichlorobenzene	ND		2.1
2-Methylphenol	ND		2.1
4-Methylphenol	ND		2.1
N-Nitrosodi-n-propylamine	ND		2.1
Hexachloroethane	ND		2.1
Nitrobenzene	ND		2.1
Isophorone	ND		2.1
2-Nitrophenol	ND		2.1
2,4-Dimethylphenol	ND		2.1
Bis(2-chloroethoxy)methane	ND		5.2
2,4-Dichlorophenol	ND		5.2
1,2,4-Trichlorobenzene	ND		2.1
Naphthalene	ND		2.1
4-Chloroaniline	ND		2.1
Hexachlorobutadiene	ND		2.1
4-Chloro-3-methylphenol	ND		5.2
2-Methylnaphthalene	ND		2.1
Hexachlorocyclopentadiene	ND		5.2
2,4,6-Trichlorophenol	ND		2.1
2,4,5-Trichlorophenol	ND		2.1
2-Chloronaphthalene	ND		2.1
2-Nitroaniline	ND		10
Dimethyl phthalate	ND		5.2
Acenaphthylene	ND		2.1
3-Nitroaniline	ND		5.2
Acenaphthene	ND		2.1
2,4-Dinitrophenol	ND		10
4-Nitrophenol	ND		10
Dibenzofuran	ND		2.1
2,4-Dinitrotoluene	ND		2.1
2.6-Dinitrotoluene	ND		5.2
Diethyl phthalate	ND		5.2
4-Chlorophenyl phenyl ether	ND		5.2
Fluorene	ND		2.1
4-Nitroaniline	ND		10
2-Methyl-4,6-dinitrophenol	ND		10
N-Nitrosodiphenylamine	ND		2.1
4-Bromophenyl phenyl ether	ND		5.2

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-8-DUP

Lab Sample ID: 720-18564-8FD Date Sampled: 03/17/2009 1245 Client Matrix: Water 03/17/2009 1545 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-

Dilution: 1.0

Initial Weight/Volume: 970 mL Date Analyzed: 03/25/2009 1757 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.1
Pentachlorophenol	ND		10
Phenanthrene	ND		2.1
Anthracene	ND		2.1
Di-n-butyl phthalate	ND		5.2
Fluoranthene	ND		2.1
Pyrene	ND		2.1
Butyl benzyl phthalate	ND		5.2
3,3'-Dichlorobenzidine	ND		5.2
Benzo[a]anthracene	ND		5.2
Bis(2-ethylhexyl) phthalate	ND		10
Chrysene	ND		2.1
Di-n-octyl phthalate	ND		21
Benzo[b]fluoranthene	ND		2.1
Benzo[a]pyrene	ND		2.1
Benzo[k]fluoranthene	ND		2.1
Indeno[1,2,3-cd]pyrene	ND		2.1
Benzo[g,h,i]perylene	ND		2.1
Benzoic acid	ND		10
Azobenzene	ND		2.1
Dibenz(a,h)anthracene	ND		2.1
Surrogate	%Rec		Acceptance Limits
Nitrobenzene-d5	69		6 - 98
2-Fluorobiphenyl	64		6 - 103
Terphenyl-d14	75		36 - 106
2-Fluorophenol	40		1 - 66
Phenol-d5	27		1 - 47
2,4,6-Tribromophenol	74		22 - 124

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: FB-031709

 Lab Sample ID:
 720-18564-9FB
 Date Sampled:
 03/17/2009 0830

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

Method: 8270C Analysis Batch: 720-48005 Instrument ID: Sat 2K

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-

Dilution: 1.0 Initial Weight/Volume: 970 mL

Date Analyzed: 03/25/2009 1831 Final Weight/Volume: 1 mL

Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Phenol ND 2.1	Analyte	Result (ug/L)	Qualifier	RL
2-Chicrophenol ND 2.1 1,3-Dichlorobenzene ND 2.1 1,4-Dichlorobenzene ND 2.1 Benzyl alcohol ND 5.2 1,2-Dichlorobenzene ND 2.1 2-Methylphenol ND 2.1 4-Methylphenol ND 2.1 N-Nitrosodi-n-propylamine ND 2.1 Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Isophorone ND 2.1 S-Nitrophenol ND 2.1 2,-Hirophenol ND 2.1 2,-Uniderlylphenol ND 2.1 2,4-Dichlorophenol ND 2.1 3,2-Tirchlorophenol ND 5.2 2,4-Dichlorophenol ND 2.1 4-Chloroa-millene ND 2.1 4-Chloroa-menylphenol ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaptrhitalene	Phenol	ND		2.1
1,3-Dichlorobenzene ND 2,1 1,4-Dichlorobenzene ND 5,2 Enenzyl alcohol ND 5,2 1,2-Dichlorobenzene ND 2,1 2-Methylphenol ND 2,1 4-Methylphenol ND 2,1 N-Nitrosodi-n-propylamine ND 2,1 Hexachloroethane ND 2,1 Nitrobenzene ND 2,1 Isophorone ND 2,1 2-Nitrophenol ND 2,1 2-Nitrophenol ND 2,1 2-Choliorophenol ND 2,1 18(2-chloroethoxy)methane ND 5,2 2,4-Dimethylphenol ND 5,2 2,4-Dindrophenol ND 5,2 2,4-Dindrophenol ND 2,1 4-Chloroanline ND 2,1 4-Chloroanline ND 2,1 4-Chloroan-methylphenol ND 2,1 4-Chloroan-methylphenol ND 2,1 4-Kaphthylaphthalene	Bis(2-chloroethyl)ether	ND		
1,4-Dichlorobenzene ND 5.2 Benzyl alcohol ND 5.2 1,2-Dichlorobenzene ND 2.1 2-Methylphenol ND 2.1 4-Methylphenol ND 2.1 N-Nitrosodi-n-propylamine ND 2.1 Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Isophorone ND 2.1 2-Nitrophenol ND 2.1 2-Nitrophenol ND 2.1 2-Nitrophenol ND 2.1 2-Nitrophenol ND 5.2 2-4-Dindrophenol ND 5.2 1,2-4-Tichlorobenzene ND 5.2 1,2-4-Tichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloro-a-methylphenol ND 2.1 Hexachlorobutadiene ND 2.1 4-Chloro-a-methylphenol ND 2.1 Hexachlorocyclopentaliene ND 5.2 2-Methylnaphthalene	2-Chlorophenol	ND		2.1
Benzyl alcohol ND 5.2 1,2-Dichlorobenzene ND 2.1 2-Methylphenol ND 2.1 4-Methylphenol ND 2.1 4-Methylphenol ND 2.1 Hexachloroethane ND 2.1 Hirobenzene ND 2.1 Isophorone ND 2.1 2-Nitrophenol ND 2.1 2-Horityphenol ND 2.1 Isi2-c-hloroethoxylmethane ND 2.1 2,4-Dichlorophenol ND 5.2 2,4-Dichlorophenol ND 5.2 2,4-Dichlorophenol ND 2.1 Naphthalene ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Sacklorocyclopentadiene ND 2.1 4-Kexaphthylphenol ND 2.1 4-Kexaphthylphenol ND 2.1 2,4-5-Trichlorophenol <td>1,3-Dichlorobenzene</td> <td>ND</td> <td></td> <td>2.1</td>	1,3-Dichlorobenzene	ND		2.1
1.2-Dichlorobenzene ND 2.1 2-Methylphenol ND 2.1 4-Methylphenol ND 2.1 N-Nitrosodi-n-propylamine ND 2.1 Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Isophorone ND 2.1 2-Nitrophenol ND 2.1 2-Nitrophenol ND 2.1 2-A-Dimethylphenol ND 2.1 Bis(2-chloroethoxy)methane ND 2.1 2,4-Dichlorophenol ND 5.2 1,2,4-Trichlorophenol ND 2.1 4-Chloro-alline ND 5.2 2-Methylnaphthalene ND 2.1 4-Exachloropthadiene ND 2.1 2,4.6-Trichlorophenol ND 2.1 2-Nitroaniline <td< td=""><td>1,4-Dichlorobenzene</td><td>ND</td><td></td><td>2.1</td></td<>	1,4-Dichlorobenzene	ND		2.1
2-Methylphenol ND 2.1 4-Methylphenol ND 2.1 N-Nitrosodi-n-propylamine ND 2.1 Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Isophorone ND 2.1 2-Nitrophenol ND 2.1 2-H-Dimethylphenol ND 2.1 Bis(2-chloroethoxy)methane ND 2.2 2,4-Dichlorophenol ND 5.2 2,4-Dichlorophenol ND 5.2 2,4-Dichlorophenol ND 2.1 Naphthalene ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Kaschlorocyclopentadiene ND 2.1 4-Kaschlorocyclopentadiene ND 2.1 4-Kaschlorophenol ND 2.1 2,4-5-Trichlorophenol ND 2.1 2-Kis-Trichlorophenol ND 2.1	Benzyl alcohol	ND		5.2
4-Methylphenol ND 2.1 N-Nitrosodi-n-propylamine ND 2.1 Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Isophorone ND 2.1 2-Nitrophenol ND 2.1 2-Nitrophenol ND 2.1 8is/2-chloroethoxy/methane ND 2.2 2-J-ichlorophenol ND 5.2 1,2-4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 Hexachlorobutadiene ND 2.1 4-Chloroa-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Kexachlorocyclopentadiene ND 2.1 4-Kexachlorocyclopentadiene ND 2.1 4-Kexachlorocyclopentadiene ND 2.1 2,4-5-Trichlorophenol ND 2.1 2,4-5-Trichlorophenol ND 2.1 2-Nitroaniline ND 2.1	1,2-Dichlorobenzene	ND		
N-Nitrosodi-n-propylamine ND 2.1 Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Isophorone ND 2.1 2-Nitrophenol ND 2.1 2-Nitrophenol ND 2.1 Sic2-chloroethoxymethane ND 5.2 2,4-Dichlorophenol ND 5.2 1,2-4-Tirchlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 Hexachlorobutadiene ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 5.2 Abethylanghthalene ND 5.2 Abethylanghthalene ND 2.1 Hexachlorocyclopentadiene ND 2.1 4-S-Tirchlorophenol ND 2.1 4-S-Tirchlorophenol ND 2.1 2-A,5-Tirchlorophenol ND 2.1 2-Nitroaniline ND 2.1 2-Nitro	2-Methylphenol	ND		
Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Isophorone ND 2.1 2-Nitrophenol ND 2.1 2,4-Dimethylphenol ND 2.1 8is(2-chloroethoxy)methane ND 5.2 2,4-Dichlorophenol ND 5.2 1,2,4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroanilline ND 2.1 4-Chloroanilline ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 5.2 4-Methylnaphthalene ND 5.2 4-Methylnaphthalene ND 5.2 4-A,6-Trichlorophenol ND 2.1 2,4,6-Trichlorophenol ND 2.1 2-A,6-Trichlorophenol ND 2.1 2-Nitroaniline ND 2.1 2-Chloronaphthalene ND 5.2 Acenaphthylene ND 5.2 Acenaphthylen	4-Methylphenol	ND		2.1
Hexachloroethane ND 2.1 Nitrobenzene ND 2.1 Sophorone ND 2.1 Sophoroethoxy)methane ND 5.2 2,4-Dichlorophenol ND 5.2 2,4-Trichlorophenol ND 2.1 Sophoroachiline ND 2.1 Hexachlorobutadiene ND 2.1 Hexachlorobutadiene ND 2.1 Hexachlorobutadiene ND 2.1 Hexachlorobutadiene ND 2.1 Hexachlorocyclopentadiene ND 5.2 Sophoroachiline ND 5.2 Sop	N-Nitrosodi-n-propylamine	ND		2.1
Sophorone		ND		2.1
2-Nitrophenol ND 2.1 2,4-Dimethylphenol ND 2.1 Bis(2-chloroethoxy)methane ND 5.2 2,4-Dichlorophenol ND 5.2 1,2,4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 5.2 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 5.2 2,4,5-Trichlorophenol ND 2.1 2-Netrophyl phthalate ND 2.1 2-Nitroaniline ND 5.2 Acenaphthylene ND 5.2 Acenaphthylene ND 10 <tr< td=""><td>Nitrobenzene</td><td>ND</td><td></td><td>2.1</td></tr<>	Nitrobenzene	ND		2.1
2,4-Dimethylphenol ND 5.2 Bis(2-chloroethoxy)methane ND 5.2 2,4-Dichlorophenol ND 5.2 1,2,4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 2.1 2-Metrylnaphthalene ND 2.1 Hexachlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Khiroraniline ND 2.1 2-Khiroraniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 2.1 4,4-Dinitrobluene ND 5.2 Dietnyl phthalate ND 5.2 4-Chlorophenyl phenyl ether <t< td=""><td>Isophorone</td><td>ND</td><td></td><td>2.1</td></t<>	Isophorone	ND		2.1
Bis(2-chloroethoxy)methane ND 5.2 2,4-Dichlorophenol ND 5.2 1,2,4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 4-Chloro-3-methylphenol ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-A,5-Trichlorophenol ND 2.1 2-A,5-Trichlorophenol ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 10 2-Nitroaniline ND 5.2 Acenaphthene ND 5.2 Acenaphthene ND 5.2 2,4-Dinitrophenol ND 10 Valitrophenol ND 2.1 2,4-Dinitrotoluene ND 5.2 Diet	2-Nitrophenol	ND		2.1
Bis(2-chloroethoxy)methane ND 5.2 2,4-Dichlorophenol ND 5.2 1,2,4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 Hexachlorobutadiene ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 5.2 Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 5.2 2,4,5-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Nitroaniline ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 2.1 Acenaphthylene ND 5.2 Acenaphthene ND 5.2 Acenaphthene ND 10 2,4-Dinitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 5.2 Diethyl phthalat		ND		2.1
1,2,4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 Hexachlorobutadiene ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-A,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 10 Acenaphthylene ND 5.2 Acenaphthene ND 2.1 2-Acenaphthene ND 2.1 2-A-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl et		ND		5.2
1,2,4-Trichlorobenzene ND 2.1 Naphthalene ND 2.1 4-Chloroaniline ND 2.1 Hexachlorobutadiene ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 5.2 Acenaphthene ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 4-Nitrophenol ND 2.1 2,4-Dinitrotoluene ND 2.1 2,4-Dinitrotoluene ND 5.2 Diethyl phthalate	2,4-Dichlorophenol	ND		5.2
Naphthalene ND 2.1 4-Chloraoniline ND 2.1 Hexachlorobutadiene ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 2.1 2-4,6-Trichlorophenol ND 2.1 2-4,5-Trichlorophenol ND 2.1 2-A,5-Trichlorophenol ND 2.1 2-A,5-Trichlorophenol ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Vibersoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether </td <td></td> <td>ND</td> <td></td> <td>2.1</td>		ND		2.1
Hexachlorobutadiene ND 2.1 4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 10 Acenaphthylene ND 5.2 Acenaphthylene ND 5.2 Acenaphthylene ND 5.2 Acenaphthene ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 4-Nitrophenol ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 5.2 4-Nitroaniline <td< td=""><td>Naphthalene</td><td>ND</td><td></td><td>2.1</td></td<>	Naphthalene	ND		2.1
4-Chloro-3-methylphenol ND 5.2 2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 4-Nitrophenol ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 5.2 4-Nitroaniline ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 </td <td>4-Chloroaniline</td> <td>ND</td> <td></td> <td>2.1</td>	4-Chloroaniline	ND		2.1
2-Methylnaphthalene ND 2.1 Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 5.2 4-Caphintrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 10 2,4-Dinitrobluene ND 2.1 2,4-Dinitrobluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 5.2 4-Nitroaniline ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1<	Hexachlorobutadiene	ND		2.1
Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 5.2 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 10 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 5.2 4-Nitroaniline ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	4-Chloro-3-methylphenol	ND		5.2
Hexachlorocyclopentadiene ND 5.2 2,4,6-Trichlorophenol ND 2.1 2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 5.2 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 10 2,4-Dinitrotoluene ND 2.1 2,4-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 5.2 4-Nitroaniline ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	2-Methylnaphthalene	ND		2.1
2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 5.2 4-Nitroaniline ND 2.1 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 10		ND		5.2
2,4,5-Trichlorophenol ND 2.1 2-Chloronaphthalene ND 2.1 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 2.1 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 10	2,4,6-Trichlorophenol	ND		2.1
2-Nitroaniline ND 10 Dimethyl phthalate ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	2,4,5-Trichlorophenol	ND		2.1
Dimethyl phthalate ND 5.2 Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	2-Chloronaphthalene	ND		2.1
Acenaphthylene ND 2.1 3-Nitroaniline ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 5.2 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	2-Nitroaniline	ND		10
3-Nitroaniline ND 5.2 Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	Dimethyl phthalate	ND		5.2
Acenaphthene ND 2.1 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	Acenaphthylene	ND		2.1
2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	3-Nitroaniline	ND		5.2
4-Nitrophenol ND 10 Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	Acenaphthene	ND		2.1
Dibenzofuran ND 2.1 2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	2,4-Dinitrophenol	ND		10
2,4-Dinitrotoluene ND 2.1 2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	4-Nitrophenol	ND		10
2,6-Dinitrotoluene ND 5.2 Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	Dibenzofuran	ND		2.1
Diethyl phthalate ND 5.2 4-Chlorophenyl phenyl ether ND 5.2 Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1	2,4-Dinitrotoluene	ND		2.1
4-Chlorophenyl phenyl etherND5.2FluoreneND2.14-NitroanilineND102-Methyl-4,6-dinitrophenolND10N-NitrosodiphenylamineND2.1	2,6-Dinitrotoluene	ND		5.2
4-Chlorophenyl phenyl etherND5.2FluoreneND2.14-NitroanilineND102-Methyl-4,6-dinitrophenolND10N-NitrosodiphenylamineND2.1	Diethyl phthalate	ND		5.2
Fluorene ND 2.1 4-Nitroaniline ND 10 2-Methyl-4,6-dinitrophenol ND 10 N-Nitrosodiphenylamine ND 2.1		ND		5.2
2-Methyl-4,6-dinitrophenolND10N-NitrosodiphenylamineND2.1		ND		2.1
2-Methyl-4,6-dinitrophenolND10N-NitrosodiphenylamineND2.1	4-Nitroaniline	ND		
N-Nitrosodiphenylamine ND 2.1				10

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: FB-031709

Lab Sample ID: 720-18564-9FB Date Sampled: 03/17/2009 0830 Client Matrix: Water 03/17/2009 1545 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Dilution: 1.0

Initial Weight/Volume: 970 mL Date Analyzed: 03/25/2009 1831 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.1
Pentachlorophenol	ND		10
Phenanthrene	ND		2.1
Anthracene	ND		2.1
Di-n-butyl phthalate	ND		5.2
Fluoranthene	ND		2.1
Pyrene	ND		2.1
Butyl benzyl phthalate	ND		5.2
3,3'-Dichlorobenzidine	ND		5.2
Benzo[a]anthracene	ND		5.2
Bis(2-ethylhexyl) phthalate	ND		10
Chrysene	ND		2.1
Di-n-octyl phthalate	ND		21
Benzo[b]fluoranthene	ND		2.1
Benzo[a]pyrene	ND		2.1
Benzo[k]fluoranthene	ND		2.1
Indeno[1,2,3-cd]pyrene	ND		2.1
Benzo[g,h,i]perylene	ND		2.1
Benzoic acid	ND		10
Azobenzene	ND		2.1
Dibenz(a,h)anthracene	ND		2.1
Surrogate	%Rec	Accep	tance Limits
Nitrobenzene-d5	60	6 - 9	8
2-Fluorobiphenyl	56	6 - 1	03
Terphenyl-d14	75	36 -	106
2-Fluorophenol	37	1 - 6	6
Phenol-d5	25	1 - 4	7
2,4,6-Tribromophenol	66	22 -	124

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-5

 Lab Sample ID:
 720-18564-1
 Date Sampled:
 03/17/2009 0935

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8015B Diesel Range Organics (DRO) (GC)

Method: 8015B Analysis Batch: 720-47867 Instrument ID: HP DRO5

Preparation: 3510C Prep Batch: 720-47825 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 500 mL Date Analyzed: 03/25/2009 0959 Final Weight/Volume: 2 mL

Date Prepared: 03/19/2009 1142 Injection Volume:

Column ID: PRIMARY

Analyte Result (ug/L) Qualifier RL

Diesel Range Organics [C10-C28] ND 50

Motor Oil Range Organics [C24-C36] ND 300

Surrogate%RecAcceptance Limitsp-Terphenyl9549 - 120

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-9

 Lab Sample ID:
 720-18564-2
 Date Sampled:
 03/17/2009 0942

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8015B Diesel Range Organics (DRO) (GC)

Method: 8015B Analysis Batch: 720-47867 Instrument ID: HP DRO5

Preparation: 3510C Prep Batch: 720-47825 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 500 mL Date Analyzed: 03/25/2009 1026 Final Weight/Volume: 2 mL

Date Prepared: 03/19/2009 1142 Injection Volume:

Column ID: PRIMARY

Analyte Result (ug/L) Qualifier RL

Diesel Range Organics [C10-C28] ND 50

Motor Oil Range Organics [C24-C36] ND 300

Surrogate%RecAcceptance Limitsp-Terphenyl9449 - 120

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-7

 Lab Sample ID:
 720-18564-3
 Date Sampled:
 03/17/2009 1115

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8015B Diesel Range Organics (DRO) (GC)

Method: 8015B Analysis Batch: 720-47867 Instrument ID: HP DRO5

Preparation: 3510C Prep Batch: 720-47825 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 500 mL Date Analyzed: 03/25/2009 0932 Final Weight/Volume: 2 mL

Date Prepared: 03/19/2009 1142 Injection Volume:

Column ID: PRIMARY

Analyte Result (ug/L) Qualifier RL

Diesel Range Organics [C10-C28] ND 50

Motor Oil Range Organics [C24-C36] ND 300

Surrogate%RecAcceptance Limitsp-Terphenyl9749 - 120

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-6

Lab Sample ID: 720-18564-4 Date Sampled: 03/17/2009 1055 Client Matrix: 03/17/2009 1545 Water Date Received:

8015B Diesel Range Organics (DRO) (GC)

8015B Analysis Batch: 720-47867 Method: Instrument ID: HP DRO5 Lab File ID: Preparation: 3510C Prep Batch: 720-47825 N/A

Dilution: 1.0

Initial Weight/Volume: 500 mL Date Analyzed: 03/25/2009 1053 Final Weight/Volume: 2 mL

Date Prepared: 03/19/2009 1142 Injection Volume:

Column ID: **PRIMARY**

Analyte Result (ug/L) Qualifier RLDiesel Range Organics [C10-C28] ND 50 Motor Oil Range Organics [C24-C36] ND 300

Surrogate %Rec Acceptance Limits 49 - 120 p-Terphenyl 106

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-8

Lab Sample ID: 720-18564-5 Date Sampled: 03/17/2009 1230 Client Matrix: Water Date Received: 03/17/2009 1545

8015B Diesel Range Organics (DRO) (GC)-Silica Gel Cleanup

Analysis Batch: 720-48126 Method: 8015B Instrument ID: HP DRO5 Preparation: 3510C SGC Prep Batch: 720-48051 Lab File ID: N/A

Dilution: 1.0

Initial Weight/Volume: 472 mL Date Analyzed: 03/28/2009 0423 Final Weight/Volume: 2 mL

Date Prepared: 03/26/2009 1703 Injection Volume:

Column ID: **PRIMARY**

Analyte Result (ug/L) Qualifier RLMotor Oil Range Organics [C24-C36] ND 320 Diesel Range Organics [C9-C24] ND 53

%Rec Acceptance Limits Surrogate 0 - 5 Capric Acid (Surr) 0 p-Terphenyl 58 31 - 120

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-2

 Lab Sample ID:
 720-18564-6
 Date Sampled:
 03/17/2009 1212

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8015B Diesel Range Organics (DRO) (GC)

Method: 8015B Analysis Batch: 720-47867 Instrument ID: HP DRO5

Preparation: 3510C Prep Batch: 720-47825 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 500 mL Date Analyzed: 03/25/2009 1158 Final Weight/Volume: 2 mL

Date Prepared: 03/19/2009 1142 Injection Volume:

Column ID: PRIMARY

Analyte Result (ug/L) Qualifier RL

Diesel Range Organics [C10-C28] ND 50

Motor Oil Range Organics [C24-C36] ND 300

Surrogate%RecAcceptance Limitsp-Terphenyl9349 - 120

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-10

 Lab Sample ID:
 720-18564-7
 Date Sampled:
 03/17/2009 1449

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8015B Diesel Range Organics (DRO) (GC)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-48126 Instrument ID: HP DRO5

Preparation: 3510C SCC Prep Batch: 720-48051 Lab File ID: N/A

Preparation: 3510C SGC Prep Batch: 720-48051 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 446 mL Date Analyzed: 03/28/2009 0450 Final Weight/Volume: 2 mL

Date Prepared: 03/26/2009 1703 Injection Volume:

Column ID: PRIMARY

Analyte	Result (ug/L)	Qualifier	RL
Motor Oil Range Organics [C24-C36]	ND		340
Diesel Range Organics [C9-C24]	ND		56

 Surrogate
 %Rec
 Acceptance Limits

 Capric Acid (Surr)
 1
 0 - 5

 p-Terphenyl
 54
 31 - 120

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: MW-8-DUP

 Lab Sample ID:
 720-18564-8FD
 Date Sampled:
 03/17/2009 1245

 Client Matrix:
 Water
 Date Received:
 03/17/2009 1545

8015B Diesel Range Organics (DRO) (GC)-Silica Gel Cleanup

Method: 8015B Analysis Batch: 720-48126 Instrument ID: HP DRO5

Preparation: 3510C SCC Prep Batch: 720-48051 Lab File ID: N/A

Preparation: 3510C SGC Prep Batch: 720-48051 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 474 mL Date Analyzed: 03/28/2009 0517 Final Weight/Volume: 2 mL

Date Prepared: 03/26/2009 1703 Injection Volume:

Column ID: PRIMARY

Analyte	Result (ug/L)	Qualifier	RL
Motor Oil Range Organics [C24-C36]	ND		320
Diesel Range Organics [C9-C24]	ND		53

Surrogate	%Rec	Acceptance Limits
Capric Acid (Surr)	0	0 - 5
p-Terphenyl	54	31 - 120

Client: LFR, Inc. Job Number: 720-18564-1

Client Sample ID: FB-031709

Lab Sample ID: 720-18564-9FB Date Sampled: 03/17/2009 0830 Client Matrix: Water Date Received: 03/17/2009 1545

8015B Diesel Range Organics (DRO) (GC)

8015B Analysis Batch: 720-47867 Method: Instrument ID: HP DRO5 Lab File ID: Preparation: 3510C Prep Batch: 720-47825 N/A

Dilution: 1.0 Initial Weight/Volume:

500 mL Date Analyzed: 03/25/2009 1319 Final Weight/Volume: 2 mL

Date Prepared: 03/19/2009 1148 Injection Volume:

Column ID: **PRIMARY**

Analyte Result (ug/L) Qualifier RLDiesel Range Organics [C10-C28] ND 50 Motor Oil Range Organics [C24-C36] ND 300

Surrogate %Rec Acceptance Limits 49 - 120 p-Terphenyl 94

DATA REPORTING QUALIFIERS

Client: LFR, Inc. Job Number: 720-18564-1

Lab Section	Qualifier	Description	
GC/MS VOA			
	F	MS or MSD exceeds the control limits	

Client: LFR, Inc. Job Number: 720-18564-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Analysis Batch:720-4	7947				
LCS 720-47947/2	Lab Control Spike	T	Water	8260B/CA_LUFT	
LCSD 720-47947/1	Lab Control Spike Duplicate	Т	Water	8260B/CA_LUFT	
MB 720-47947/3	Method Blank	Т	Water	8260B/CA_LUFT	
720-18564-1	MW-5	Т	Water	8260B/CA_LUFT	
720-18564-1MS	Matrix Spike	Т	Water	8260B/CA_LUFT	
720-18564-1MSD	Matrix Spike Duplicate	T	Water	8260B/CA_LUFT	
720-18564-2	MW-9	Т	Water	8260B/CA_LUFT	
720-18564-3	MW-7	T	Water	8260B/CA LUFT	
720-18564-4	MW-6	T	Water	8260B/CA LUFT	
720-18564-5	MW-8	T	Water	8260B/CA LUFT	
720-18564-6	MW-2	T	Water	8260B/CA LUFT	
720-18564-7	MW-10	T	Water	8260B/CA_LUFT	
720-18564-8FD	MW-8-DUP	T	Water	8260B/CA LUFT	
720-18564-9FB	FB-031709	T	Water	8260B/CA LUFT	

Report Basis

T = Total

Client: LFR, Inc. Job Number: 720-18564-1

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC/MS Semi VOA					
Prep Batch: 720-47916	1				
LCS 720-47916/2-A	Lab Control Spike	T	Water	3510C	
LCSD 720-47916/3-A	Lab Control Spike Duplicate	T	Water	3510C	
MB 720-47916/1-A	Method Blank	T	Water	3510C	
720-18564-1	MW-5	T	Water	3510C	
720-18564-2	MW-9	T	Water	3510C	
720-18564-3	MW-7	T	Water	3510C	
720-18564-4	MW-6	T	Water	3510C	
720-18564-5	MW-8	T	Water	3510C	
720-18564-6	MW-2	T	Water	3510C	
720-18564-7	MW-10	T	Water	3510C	
720-18564-8FD	MW-8-DUP	T	Water	3510C	
720-18564-9FB	FB-031709	T	Water	3510C	
Analysis Batch:720-48	005				
LCS 720-47916/2-A	Lab Control Spike	T	Water	8270C	720-47916
LCSD 720-47916/3-A	Lab Control Spike Duplicate	T	Water	8270C	720-47916
MB 720-47916/1-A	Method Blank	T	Water	8270C	720-47916
720-18564-1	MW-5	T	Water	8270C	720-47916
720-18564-2	MW-9	T	Water	8270C	720-47916
720-18564-3	MW-7	T	Water	8270C	720-47916
720-18564-4	MW-6	Т	Water	8270C	720-47916
720-18564-5	MW-8	T	Water	8270C	720-47916
720-18564-6	MW-2	T	Water	8270C	720-47916
720-18564-7	MW-10	T	Water	8270C	720-47916
720-18564-8FD	MW-8-DUP	Т	Water	8270C	720-47916
720-18564-9FB	FB-031709	Т	Water	8270C	720-47916

Report Basis

T = Total

Client: LFR, Inc. Job Number: 720-18564-1

QC Association Summary

•		Danari			
Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC Semi VOA					
Prep Batch: 720-47825					
LCS 720-47825/2-A	Lab Control Spike	T	Water	3510C	
LCSD 720-47825/3-A	Lab Control Spike Duplicate	T	Water	3510C	
MB 720-47825/1-A	Method Blank	T	Water	3510C	
720-18564-1	MW-5	Т	Water	3510C	
720-18564-2	MW-9	T	Water	3510C	
720-18564-3	MW-7	T	Water	3510C	
720-18564-4	MW-6	T	Water	3510C	
720-18564-6	MW-2	Т	Water	3510C	
720-18564-9FB	FB-031709	T	Water	3510C	
Analysis Batch:720-47	867				
LCS 720-47825/2-A	Lab Control Spike	Т	Water	8015B	720-47825
LCSD 720-47825/3-A	Lab Control Spike Duplicate	Т	Water	8015B	720-47825
MB 720-47825/1-A	Method Blank	Т	Water	8015B	720-47825
720-18564-1	MW-5	Т	Water	8015B	720-47825
720-18564-2	MW-9	Т	Water	8015B	720-47825
720-18564-3	MW-7	Т	Water	8015B	720-47825
720-18564-4	MW-6	Т	Water	8015B	720-47825
720-18564-6	MW-2	Т	Water	8015B	720-47825
720-18564-9FB	FB-031709	Т	Water	8015B	720-47825
Prep Batch: 720-48051					
LCS 720-48051/2-A	Lab Control Spike	Α	Water	3510C SGC	
LCSD 720-48051/3-A	Lab Control Spike Duplicate	Α	Water	3510C SGC	
MB 720-48051/1-A	Method Blank	Α	Water	3510C SGC	
720-18564-5	MW-8	Α	Water	3510C SGC	
720-18564-7	MW-10	Α	Water	3510C SGC	
720-18564-8FD	MW-8-DUP	Α	Water	3510C SGC	
Analysis Batch:720-48	126				
LCS 720-48051/2-A	Lab Control Spike	Α	Water	8015B	720-48051
LCSD 720-48051/3-A	Lab Control Spike Duplicate	A	Water	8015B	720-48051
MB 720-48051/1-A	Method Blank	A	Water	8015B	720-48051
720-18564-5	MW-8	A	Water	8015B	720-48051
720-18564-7	MW-10	A	Water	8015B	720-48051
720-18564-8FD	MW-8-DUP	A	Water	8015B	720-48051
20 10001 01 D		, ,		30100	120 40001

Report Basis

A = Silica Gel Cleanup

T = Total

Client: LFR, Inc. Job Number: 720-18564-1

Method Blank - Batch: 720-47947 Method: 8260B/CA_LUFTMS

Preparation: 5030B

Lab Sample ID: MB 720-47947/3 Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Client Matrix: Water Prep Batch: N/A Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL

Date Analyzed: 03/23/2009 0955 Final Weight/Volume: 10 mL Date Prepared: 03/23/2009 0955

Analyte	Result	Qual	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	% Rec	Acceptance Limit	S
Toluene-d8 (Surr)	93	78 - 112	
1,2-Dichloroethane-d4 (Surr)	87	67 - 126	

Client: LFR, Inc. Job Number: 720-18564-1

Lab Control Spike/ Method: 8260B/CA_LUFTMS

Lab Control Spike Duplicate Recovery Report - Batch: 720-47947 Preparation: 5030B

LCS Lab Sample ID: LCS 720-47947/2 Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Client Matrix: Water Prep Batch: N/A Lab File ID: e:\data\2009\200903\03230

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL Date Analyzed: 03/23/2009 1028 Final Weight/Volume: 10 mL

Date Prepared: 03/23/2009 1028

LCSD Lab Sample ID: LCSD 720-47947/1 Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Client Matrix: Water Prep Batch: N/A Lab File ID: e:\data\2009\200903\03230\$

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL

Date Analyzed: 03/23/2009 1052 Final Weight/Volume: 10 mL

Date Prepared:

03/23/2009 1052

% Rec. LCS **LCSD RPD** RPD Limit LCS Qual LCSD Qual Analyte Limit 74 - 112 Benzene 93 90 4 20 Gasoline Range Organics (GRO)-C5-C12 69 66 42 - 80 4 20 Toluene 76 78 65 - 98 2 20 **MTBE** 88 69 - 104 20 80 8 Surrogate LCS % Rec LCSD % Rec Acceptance Limits 92 Toluene-d8 (Surr) 93 78 - 112 1,2-Dichloroethane-d4 (Surr) 90 80 67 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Matrix Spike/

Matrix Spike Duplicate Recovery Report - Batch: 720-47947

Method: 8260B/CA_LUFTMS

Preparation: 5030B

MS Lab Sample ID: 720-18564-1

Client Matrix: Water

Dilution: 1.0

Date Analyzed: 03/23/2009 1517 Date Prepared: 03/23/2009 1517 Analysis Batch: 720-47947

Prep Batch: N/A

Instrument ID: Varian 3900A

Lab File ID: e:\data\2009\200903\0323

Initial Weight/Volume: 10 mL Final Weight/Volume: 10 mL

MSD Lab Sample ID: 720-18564-1

Client Matrix: Water Dilution: 1.0

Date Analyzed: 03/23/2009 1539 Date Prepared: 03/23/2009 1539 Analysis Batch: 720-47947 Instrument ID: Varian 3900A

Prep Batch: N/A Lab File ID: e:\data\2009\200903\03230

Initial Weight/Volume: 10 mL Final Weight/Volume: 10 mL

	<u>%</u>	<u> 6 Rec.</u>				
Analyte	MS	MSD	Limit	RPD	RPD Limit	MS Qual MSD Qual
Benzene	87	89	58 - 134	2	20	
Gasoline Range Organics (GRO)-C5-C12	59	60	43 - 95	1	20	
Toluene	70	73	72 - 129	5	20	F
MTBE	89	84	22 - 185	5	20	
Surrogate		MS % Rec	MSD ^c	% Rec	Acce	eptance Limits
Toluene-d8 (Surr)		93	90		78	8 - 112
1,2-Dichloroethane-d4 (Surr)		91	86		6	7 - 126

Client: LFR, Inc. Job Number: 720-18564-1

Method Blank - Batch: 720-47916 Method: 8270C Preparation: 3510C

Lab Sample ID: MB 720-47916/1-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\mb

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL
Date Analyzed: 03/25/2009 1149

Date Prepared: 03/23/2009 1257

Injection Volume: 1.0 uL

Qual RL Analyte Result Phenol ND 2.0 Bis(2-chloroethyl)ether ND 2.0 2-Chlorophenol ND 2.0 1,3-Dichlorobenzene ND 2.0 1,4-Dichlorobenzene ND 2.0 Benzyl alcohol 5.0 ND 1,2-Dichlorobenzene ND 2.0 2-Methylphenol ND 2.0 4-Methylphenol 2.0 ND N-Nitrosodi-n-propylamine ND 2.0 Hexachloroethane ND 2.0 Nitrobenzene 2.0 ND 2.0 Isophorone ND 2-Nitrophenol ND 2.0 2.0 2,4-Dimethylphenol ND Bis(2-chloroethoxy)methane 5.0 ND 5.0 2,4-Dichlorophenol ND 1,2,4-Trichlorobenzene ND 2.0 Naphthalene ND 2.0 4-Chloroaniline ND 2.0 Hexachlorobutadiene ND 2.0 4-Chloro-3-methylphenol ND 5.0 2-Methylnaphthalene ND 2.0 Hexachlorocyclopentadiene ND 5.0 2,4,6-Trichlorophenol ND 2.0 2,4,5-Trichlorophenol ND 2.0 2-Chloronaphthalene ND 2.0 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.0 Acenaphthylene ND 2.0 3-Nitroaniline ND 5.0 Acenaphthene ND 2.0 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.0 2.4-Dinitrotoluene 2.0 ND 2,6-Dinitrotoluene ND 5.0 Diethyl phthalate ND 5.0 4-Chlorophenyl phenyl ether 5.0 ND Fluorene ND 2.0 4-Nitroaniline ND 10

Client: LFR, Inc. Job Number: 720-18564-1

Method Blank - Batch: 720-47916

Method: 8270C Preparation: 3510C

Lab Sample ID: MB 720-47916/1-A

Client Matrix: Water

Dilution: 1.0

Date Analyzed: 03/25/2009 1149 Date Prepared: 03/23/2009 1257 Analysis Batch: 720-48005 Prep Batch: 720-47916

Units: ug/L

Instrument ID: Sat 2K1

Lab File ID: d:\data\200903\032509\mb

Initial Weight/Volume: 1000 mL Final Weight/Volume: 1 mL Injection Volume: 1.0 uL

Analyte	Result	Qual	RL
2-Methyl-4,6-dinitrophenol	ND		10
N-Nitrosodiphenylamine	ND		2.0
4-Bromophenyl phenyl ether	ND		5.0
Hexachlorobenzene	ND		2.0
Pentachlorophenol	ND		10
Phenanthrene	ND		2.0
Anthracene	ND		2.0
Di-n-butyl phthalate	ND		5.0
Fluoranthene	ND		2.0
Pyrene	ND		2.0
Butyl benzyl phthalate	ND		5.0
3,3'-Dichlorobenzidine	ND		5.0
Benzo[a]anthracene	ND		5.0
Bis(2-ethylhexyl) phthalate	ND		10
Chrysene	ND		2.0
Di-n-octyl phthalate	ND		20
Benzo[b]fluoranthene	ND		2.0
Benzo[a]pyrene	ND		2.0
Benzo[k]fluoranthene	ND		2.0
Indeno[1,2,3-cd]pyrene	ND		2.0
Benzo[g,h,i]perylene	ND		2.0
Benzoic acid	ND		10
Azobenzene	ND		2.0
Dibenz(a,h)anthracene	ND		2.0
Surrogate	% Rec	Acceptance Limits	
Nitrobenzene-d5	72	6 - 98	
2-Fluorobiphenyl	69	6 - 103	
Terphenyl-d14	84	36 - 106	
2-Fluorophenol	46	1 - 66	
Phenol-d5	32	1 - 47	
2,4,6-Tribromophenol	74	22 - 124	

Client: LFR, Inc. Job Number: 720-18564-1

Lab Control Spike/ Method: 8270C
Lab Control Spike Duplicate Recovery Report - Batch: 720-47916 Preparation: 3510C

LCS Lab Sample ID: LCS 720-47916/2-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\LC

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL Date Analyzed: 03/25/2009 1042 Final Weight/Volume: 1 mL

Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

LCSD Lab Sample ID: LCSD 720-47916/3-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\LCS

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL

 Date Analyzed:
 03/25/2009 1116
 Final Weight/Volume:
 1 mL

 Date Prepared:
 03/23/2009 1257
 Injection Volume:
 1.0 uL

	<u>0</u>	% Rec.					
Analyte	LCS	LCSD	Limit	RPD	RPD Limit	LCS Qual	LCSD Qual
Phenol	31	38	10 - 49	22	35		
Bis(2-chloroethyl)ether	58	71	22 - 113	21	35		
2-Chlorophenol	60	76	19 - 104	24	25		
1,3-Dichlorobenzene	62	70	18 - 95	13	35		
1,4-Dichlorobenzene	54	67	17 - 82	22	30		
Benzyl alcohol	60	75	21 - 94	22	35		
1,2-Dichlorobenzene	64	71	20 - 89	10	35		
2-Methylphenol	58	70	18 - 100	19	35		
4-Methylphenol	51	62	16 - 93	19	35		
N-Nitrosodi-n-propylamine	64	81	34 - 108	23	34		
Hexachloroethane	64	75	4 - 87	16	35		
Nitrobenzene	74	96	27 - 117	26	35		
Isophorone	79	95	42 - 118	19	35		
2-Nitrophenol	67	82	34 - 107	20	35		
2,4-Dimethylphenol	81	98	42 - 122	19	35		
Bis(2-chloroethoxy)methane	65	77	37 - 110	18	35		
2,4-Dichlorophenol	67	81	39 - 107	20	35		
1,2,4-Trichlorobenzene	62	74	20 - 95	17	35		
Naphthalene	65	80	22 - 99	21	35		
4-Chloroaniline	45	54	10 - 80	17	35		
Hexachlorobutadiene	52	66	3 - 101	24	35		
4-Chloro-3-methylphenol	74	88	22 - 147	17	31		
2-Methylnaphthalene	64	77	10 - 130	19	35		
Hexachlorocyclopentadiene	65	79	21 - 101	19	35		
2,4,6-Trichlorophenol	72	81	31 - 118	12	35		
2,4,5-Trichlorophenol	71	83	35 - 112	16	35		
2-Chloronaphthalene	63	79	26 - 106	22	35		
2-Nitroaniline	72	86	31 - 106	18	35		
Dimethyl phthalate	72	88	47 - 117	20	35		
Acenaphthylene	83	103	33 - 128	22	35		
3-Nitroaniline	71	85	40 - 107	18	35		
Acenaphthene	75	81	31 - 109	8	30		

Client: LFR, Inc. Job Number: 720-18564-1

Lab Control Spike/ Method: 8270C
Lab Control Spike Duplicate Recovery Report - Batch: 720-47916 Preparation: 3510C

LCS Lab Sample ID: LCS 720-47916/2-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\LC

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL Date Analyzed: 03/25/2009 1042 Final Weight/Volume: 1 mL

Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

LCSD Lab Sample ID: LCSD 720-47916/3-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\LCS

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL

 Date Analyzed:
 03/25/2009 1116
 Final Weight/Volume:
 1 mL

 Date Prepared:
 03/23/2009 1257
 Injection Volume:
 1.0 uL

	0	% Rec.					
Analyte	LCS	LCSD	Limit	RPD	RPD Limit	LCS Qual	LCSD Qual
2,4-Dinitrophenol	67	76	51 - 102	13	35		
4-Nitrophenol	42	48	26 - 63	13	35		
Dibenzofuran	68	79	42 - 97	14	35		
2,4-Dinitrotoluene	77	92	52 - 113	18	35		
2,6-Dinitrotoluene	77	83	50 - 115	8	35		
Diethyl phthalate	76	89	54 - 111	16	35		
4-Chlorophenyl phenyl ether	73	84	40 - 110	14	35		
Fluorene	71	85	37 - 104	17	35		
4-Nitroaniline	79	89	56 - 116	12	35		
2-Methyl-4,6-dinitrophenol	66	82	47 - 114	21	35		
N-Nitrosodiphenylamine	80	96	56 - 123	19	35		
4-Bromophenyl phenyl ether	76	95	52 - 111	22	35		
Hexachlorobenzene	78	93	61 - 104	18	35		
Pentachlorophenol	67	80	55 - 107	18	35		
Phenanthrene	75	91	56 - 110	19	35		
Anthracene	80	94	58 - 114	16	35		
Di-n-butyl phthalate	86	104	56 - 114	19	35		
Fluoranthene	83	96	60 - 121	15	35		
Pyrene	76	83	56 - 91	10	35		
Butyl benzyl phthalate	73	82	37 - 100	12	35		
3,3'-Dichlorobenzidine	59	68	37 - 111	15	35		
Benzo[a]anthracene	67	79	50 - 112	16	35		
Bis(2-ethylhexyl) phthalate	77	88	59 - 111	13	35		
Chrysene	73	77	56 - 94	4	35		
Di-n-octyl phthalate	71	83	47 - 118	15	35		
Benzo[b]fluoranthene	75	80	55 - 110	7	35		
Benzo[a]pyrene	66	75	59 - 103	12	35		
Benzo[k]fluoranthene	73	89	55 - 110	20	35		
Indeno[1,2,3-cd]pyrene	78	88	63 - 126	11	35		
Benzo[g,h,i]perylene	79	89	10 - 140	12	35		
Benzoic acid	28	30	7 - 46	8	35		
Azobenzene	76	87	49 - 102	13	35		

Client: LFR, Inc. Job Number: 720-18564-1

Lab Control Spike/ Method: 8270C
Lab Control Spike Duplicate Recovery Report - Batch: 720-47916 Preparation: 3510C

LCS Lab Sample ID: LCS 720-47916/2-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\LC

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL

 Date Analyzed:
 03/25/2009 1042
 Final Weight/Volume:
 1 mL

 Date Prepared:
 03/23/2009 1257
 Injection Volume:
 1.0 uL

LCSD Lab Sample ID: LCSD 720-47916/3-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\LCS

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL

 Date Analyzed:
 03/25/2009 1116
 Final Weight/Volume:
 1 mL

 Date Prepared:
 03/23/2009 1257
 Injection Volume:
 1.0 uL

% Rec. LCS **LCSD RPD** RPD Limit LCS Qual LCSD Qual Analyte Limit Dibenz(a,h)anthracene 69 82 10 - 130 17 35 Surrogate LCS % Rec LCSD % Rec Acceptance Limits Nitrobenzene-d5 70 86 6 - 98 2-Fluorobiphenyl 76 64 6 - 103 Terphenyl-d14 74 83 36 - 106 2-Fluorophenol 40 47 1 - 66 Phenol-d5 29 35 1 - 47 2,4,6-Tribromophenol 75 83 22 - 124

500 mL

49 - 120

Client: LFR, Inc. Job Number: 720-18564-1

Method Blank - Batch: 720-47825 Method: 8015B Preparation: 3510C

Lab Sample ID: MB 720-47825/1-A Analysis Batch: 720-47867 Instrument ID: HP DRO5

Client Matrix: Water Prep Batch: 720-47825 Lab File ID: N/A

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 500 mL Date Analyzed: 03/20/2009 0549 Final Weight/Volume: 2 mL

Date Prepared: 03/19/2009 1142 Injection Volume:

Column ID: PRIMARY

Analyte Result Qual RL

Diesel Range Organics [C10-C28] ND 50

Motor Oil Range Organics [C24-C36] ND 300

Surrogate % Rec Acceptance Limits p-Terphenyl 98 49 - 120

Lab Control Spike/ Method: 8015B
Lab Control Spike Duplicate Recovery Report - Batch: 720-47825 Preparation: 3510C

LCS Lab Sample ID: LCS 720-47825/2-A Analysis Batch: 720-47867 Instrument ID: HP DRO5

Client Matrix: Water Prep Batch: 720-47825 Lab File ID: N/A

Dilution: 1.0 Units: ug/L Initial Weight/Volume:

Date Analyzed: 03/20/2009 0455 Final Weight/Volume: 2 mL

Date Prepared: 03/19/2009 1142 Injection Volume:

Column ID: PRIMARY

LCSD Lab Sample ID: LCSD 720-47825/3-A Analysis Batch: 720-47867 Instrument ID: HP DRO5

Client Matrix: Water Prep Batch: 720-47825 Lab File ID: N/A

108

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 500 mL Date Analyzed: 03/20/2009 0522 Final Weight/Volume: 2 mL

Date Prepared: 03/19/2009 1142 Injection Volume:

Column ID: PRIMARY

% Rec. LCS **RPD** RPD Limit LCS Qual LCSD Qual Analyte LCSD Limit 47 - 111 Diesel Range Organics [C10-C28] 103 97 6 30 LCS % Rec LCSD % Rec Surrogate Acceptance Limits

103

Calculations are performed before rounding to avoid round-off errors in calculated results.

p-Terphenyl

Client: LFR, Inc. Job Number: 720-18564-1

Method Blank - Batch: 720-48051 Method: 8015B

Preparation: 3510C SGC Silica Gel Cleanup

Lab Sample ID: MB 720-48051/1-A

Client Matrix: Water Dilution: 1.0

Date Analyzed: 03/28/2009 0328 Date Prepared: 03/26/2009 1703 Analysis Batch: 720-48126 Prep Batch: 720-48051

Units: ug/L

Instrument ID: HP DRO5 Lab File ID: N/A

Initial Weight/Volume: 500 mL Final Weight/Volume: 2 mL

Injection Volume:

Column ID: PRIMARY

Analyte	Result	Qual RL	
Diesel Range Organics [C9-C24]	ND	50	
Motor Oil Range Organics [C24-C36]	ND	300)
Surrogate	% Rec	Acceptance Limits	
Capric Acid (Surr)	0	0 - 5	

Lab Control Spike/ Method: 8015B

Lab Control Spike Duplicate Recovery Report - Batch: 720-48051 Preparation: 3510C SGC

Silica Gel Cleanup

LCS Lab Sample ID: LCS 720-48051/2-A Analysis Batch: 720-48126 Instrument ID: HP DRO5

Client Matrix: Water Prep Batch: 720-48051 Lab File ID: N/A
Dilution: 1.0 Units: ug/L Initial Weight/Volume:

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 500 mL
Date Analyzed: 03/28/2009 0234 Final Weight/Volume: 2 mL
Date Prepared: 03/26/2009 1703 Injection Volume:

Column ID: PRIMARY

LCSD Lab Sample ID: LCSD 720-48051/3-A Analysis Batch: 720-48126 Instrument ID: HP DRO5

Client Matrix: Water Prep Batch: 720-48051 Lab File ID: N/A
Dilution: 1.0 Units: ug/L Initial Weight/Volume: 500 mL

Date Analyzed: 03/28/2009 0301 Final Weight/Volume: 2 mL

Date Prepared: 03/26/2009 1703 Injection Volume: Column ID: F

Column ID: PRIMARY

<u>% Rec.</u>

Analyte LCS **LCSD** Limit **RPD** RPD Limit LCS Qual LCSD Qual Diesel Range Organics [C9-C24] 76 81 41 - 103 30 6 LCS % Rec LCSD % Rec Surrogate Acceptance Limits 92 90 31 - 120 p-Terphenyl

720-18564 CHAIN OF CUSTODY / ANALYSES REQUEST FORM PROJECT NO .: SECTION NO .: SAMPLER'S INITIALS: 3/17/09 001-09567-07 1900 Powell Street, 12th Floor Emeryville, California 94608 (510) 652-4500 Fax: (510) 652-2246 PROJECT NAME: Hasson Radum SAMPLER (Signature): ANALYSES REMARKS SAMPLE Health fea entiringo Bret ternerneen VOCS lega executora TYPE TAT Tand Eas south *VOCs: **Metals: ☐ 8260 List ☐ CAM17 RUSH ☐ 8240 List ☐ RCRA ☐ 8010 List ☐ LUFT TIME DATE SAMPLE ID. ☐ 624 List 3/17 4:35 NW-5 MW-9 9:42 6 Z MW-7 11:15 6 MW-6 10:55 X 12:30 MW-8 5 X 6 MW-2 X 12:12 6 X X 14:49 MW-10 X MW-8-DUP 6 X TB-031704 TB 2 X 10 Page 2 RELINQUISHED BY: SAMPLE RECEIPT: RELINQUISHED BY: RELINQUISHED BY: Cooler Temp: METHOD OF SHIPMENT 1,4,2,400 ☐Intact ☐Cold (DATE) (SIGNATURE) (DATE) (SIGNATURE) LAB REPORT NO .: On Ice Ambient Cooler No: Tom Collins 15:45 (PRINTED NAME) (TIME) (TIME) (PRINTED NAME) (TIME) (PRINTED NAME) FAX COC CONFIRMATION TO: LFR Preservative Correct? Ron Goldbow Yes No NA (COMPANY) (COMPANY) (COMPANY) FAX RESULTS TO: 11 ANALYTICAL LABORATORY: RECEIVED BY: RECEIVED BY: 2 RECEIVED BY (LABORATORY): (SIGNATURE) (SIGNATURE) (DATE) (SIGNATURE) (DATE) SEND HARDCOPY TO: TBHILOCK PRINTED NAME) (PRINTED NAME) (TIME) (PRINTED NAME) (TIME) SEND EDD TO: TEST AMERICA EMV.LABEDDS.COM (COMPANY) (COMPANY) CHAIN of CUSTODY - ANALYSES FORM CDR 5/2003 Field Copy (Pink) Shipping Copy (White) File Copy (Yellow)

Login Sample Receipt Check List

Client: LFR, Inc. Job Number: 720-18564-1

Login Number: 18564 List Source: TestAmerica San Francisco

Creator: Bullock, Tracy List Number: 1

Question	T / F/ NA	Comment
Radioactivity either was not measured or, if measured, is at or below background	N/A	
The cooler's custody seal, if present, is intact.	N/A	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the sample IDs on the containers and the COC.	False	ncm
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	



ANALYTICAL REPORT

Job Number: 720-18554-1

Job Description: Hanson Radum

For: LFR, Inc. 1900 Powell St 12th Floor Emeryville, CA 94608-1827

Attention: Mr. Ron Goloubow

Approved for release Afsaneh Salimpour Project Manager I 4/30/2009 12:11 PM

Afsaneh Salimpour Project Manager I afsaneh.salimpour@testamericainc.com 04/30/2009 Revision: 1

Akanef Sal

Job Narrative 720-J18554-1

Comments

No additional comments.

Receipt

No client label on the one Trip Blank TAL-SF TB: 021709 received, no sample date collected provided used 03/16/09.

All other samples were received in good condition within temperature requirements.

GC/MS VOA

No analytical or quality issues were noted.

GC/MS Semi VOA

No analytical or quality issues were noted.

GC Semi VOA

No analytical or quality issues were noted.

Organic Prep

No analytical or quality issues were noted.

EXECUTIVE SUMMARY - Detections

Client: LFR, Inc. Job Number: 720-18554-1

Lab Sample ID Client Sample ID Reporting
Analyte Result / Qualifier Limit Units Method

No Detections

METHOD SUMMARY

Client: LFR, Inc. Job Number: 720-18554-1

Description	Lab Location	Method	Preparation Method
Matrix: Water			
Volatile Organic Compounds by GC/MS Purge and Trap	TAL SF TAL SF	SW846 8260	DB/CA_LUFTMS SW846 5030B
Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)	TAL SF	SW846 8270	OC
Liquid-Liquid Extraction (Separatory Funnel)	TAL SF		SW846 3510C

Lab References:

TAL SF = TestAmerica San Francisco

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

METHOD / ANALYST SUMMARY

Client: LFR, Inc. Job Number: 720-18554-1

Method	Analyst	Analyst ID
SW846 8260B/CA_LUFTMS	Zhao, June	JZ
SW846 8270C	Lee. Michael	ML

SAMPLE SUMMARY

Client: LFR, Inc. Job Number: 720-18554-1

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received	
720-18554-1	MW-1	Water	03/16/2009 1440	03/16/2009 1630	
720-18554-2	MW-3	Water	03/16/2009 1505	03/16/2009 1630	

Client: LFR, Inc. Job Number: 720-18554-1

Client Sample ID: MW-1

 Lab Sample ID:
 720-18554-1
 Date Sampled:
 03/16/2009 1440

 Client Matrix:
 Water
 Date Received:
 03/16/2009 1630

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA LUFTMS Analysis Batch: 720-47847 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03190

Dilution: 1.0 Initial Weight/Volume: 10 mL Date Analyzed: 03/19/2009 1832 Final Weight/Volume: 10 mL

Date Prepared: 03/19/2009 1832

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	91		78 - 112
1,2-Dichloroethane-d4 (Surr)	86		67 - 126

Client: LFR, Inc. Job Number: 720-18554-1

Client Sample ID: MW-3

 Lab Sample ID:
 720-18554-2
 Date Sampled:
 03/16/2009 1505

 Client Matrix:
 Water
 Date Received:
 03/16/2009 1630

8260B/CA_LUFTMS Volatile Organic Compounds by GC/MS

Method: 8260B/CA_LUFTMS Analysis Batch: 720-47847 Instrument ID: Varian 3900A

Preparation: 5030B Lab File ID: e:\data\2009\200903\03190

Dilution: 1.0 Initial Weight/Volume: 10 mL Date Analyzed: 03/19/2009 1855 Final Weight/Volume: 10 mL

Date Prepared: 03/19/2009 1855

Analyte	Result (ug/L)	Qualifier	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	%Rec		Acceptance Limits
Toluene-d8 (Surr)	93		78 - 112
1,2-Dichloroethane-d4 (Surr)	91		67 - 126

Client: LFR, Inc. Job Number: 720-18554-1

Client Sample ID: MW-1

 Lab Sample ID:
 720-18554-1
 Date Sampled:
 03/16/2009 1440

 Client Matrix:
 Water
 Date Received:
 03/16/2009 1630

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

Method: 8270C Analysis Batch: 720-48005 Instrument ID: Sat 2K²

Preparation: 3510C Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\720-

Dilution: 1.0 Initial Weight/Volume: 890 mL
Date Analyzed: 03/25/2009 1256 Final Weight/Volume: 1 mL
Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Result (ug/L) Qualifier RL Analyte Phenol ND 2.2 ND Bis(2-chloroethyl)ether 2.2 2-Chlorophenol 2.2 ND 1,3-Dichlorobenzene 2.2 ND 1,4-Dichlorobenzene ND 2.2 Benzvl alcohol ND 5.6 1,2-Dichlorobenzene ND 2.2 2-Methylphenol ND 2.2 4-Methylphenol ND 2.2 N-Nitrosodi-n-propylamine ND 2.2 Hexachloroethane ND 2.2 Nitrobenzene ND 2.2 Isophorone ND 2.2 2-Nitrophenol 2.2 ND 2,4-Dimethylphenol ND 2.2 Bis(2-chloroethoxy)methane ND 5.6 2,4-Dichlorophenol ND 5.6 1,2,4-Trichlorobenzene ND 2.2 Naphthalene ND 2.2 4-Chloroaniline ND 2.2 Hexachlorobutadiene ND 2.2 4-Chloro-3-methylphenol ND 5.6 2-Methylnaphthalene ND 2.2 Hexachlorocyclopentadiene ND 5.6 2,4,6-Trichlorophenol ND 2.2 2,4,5-Trichlorophenol ND 2.2 2-Chloronaphthalene ND 2.2 2-Nitroaniline ND 11 Dimethyl phthalate ND 5.6 Acenaphthylene ND 2.2 3-Nitroaniline ND 5.6 Acenaphthene ND 2.2 2,4-Dinitrophenol ND 11 4-Nitrophenol ND 11 Dibenzofuran ND 2.2 2.2 2,4-Dinitrotoluene ND 2,6-Dinitrotoluene ND 5.6 Diethyl phthalate ND 5.6 4-Chlorophenyl phenyl ether ND 5.6 Fluorene ND 2.2 4-Nitroaniline ND 11 2-Methyl-4,6-dinitrophenol ND 11 N-Nitrosodiphenylamine ND 2.2 4-Bromophenyl phenyl ether ND 5.6

Client: LFR, Inc. Job Number: 720-18554-1

Client Sample ID: MW-1

Lab Sample ID: 720-18554-1 Date Sampled: 03/16/2009 1440 Client Matrix: Water 03/16/2009 1630 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Dilution: 1.0

Initial Weight/Volume: 890 mL Date Analyzed: 03/25/2009 1256 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.2
Pentachlorophenol	ND		11
Phenanthrene	ND		2.2
Anthracene	ND		2.2
Di-n-butyl phthalate	ND		5.6
Fluoranthene	ND		2.2
Pyrene	ND		2.2
Butyl benzyl phthalate	ND		5.6
3,3'-Dichlorobenzidine	ND		5.6
Benzo[a]anthracene	ND		5.6
Bis(2-ethylhexyl) phthalate	ND		11
Chrysene	ND		2.2
Di-n-octyl phthalate	ND		22
Benzo[b]fluoranthene	ND		2.2
Benzo[a]pyrene	ND		2.2
Benzo[k]fluoranthene	ND		2.2
Indeno[1,2,3-cd]pyrene	ND		2.2
Benzo[g,h,i]perylene	ND		2.2
Benzoic acid	ND		11
Azobenzene	ND		2.2
Dibenz(a,h)anthracene	ND		2.2
Surrogate	%Rec	Ac	ceptance Limits
Nitrobenzene-d5	71	6	- 98
2-Fluorobiphenyl	70	6	- 103
Terphenyl-d14	78	3	6 - 106
2-Fluorophenol	44	1	- 66
Phenol-d5	31	1	- 47
2,4,6-Tribromophenol	81	2	2 - 124

Client: LFR, Inc. Job Number: 720-18554-1

Client Sample ID: MW-3

Lab Sample ID: 720-18554-2 Date Sampled: 03/16/2009 1505 Client Matrix: Water Date Received: 03/16/2009 1630

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID: 950 mL

Dilution: 1.0

Initial Weight/Volume: Date Analyzed: 03/25/2009 1329 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Phenol	ND		2.1
Bis(2-chloroethyl)ether	ND		2.1
2-Chlorophenol	ND		2.1
1,3-Dichlorobenzene	ND		2.1
1,4-Dichlorobenzene	ND		2.1
Benzyl alcohol	ND		5.3
1,2-Dichlorobenzene	ND		2.1
2-Methylphenol	ND		2.1
4-Methylphenol	ND		2.1
N-Nitrosodi-n-propylamine	ND		2.1
Hexachloroethane	ND		2.1
Nitrobenzene	ND		2.1
Isophorone	ND		2.1
2-Nitrophenol	ND		2.1
2,4-Dimethylphenol	ND		2.1
Bis(2-chloroethoxy)methane	ND		5.3
2,4-Dichlorophenol	ND		5.3
1,2,4-Trichlorobenzene	ND		2.1
Naphthalene	ND		2.1
4-Chloroaniline	ND		2.1
Hexachlorobutadiene	ND		2.1
4-Chloro-3-methylphenol	ND		5.3
2-Methylnaphthalene	ND		2.1
Hexachlorocyclopentadiene	ND		5.3
2,4,6-Trichlorophenol	ND		2.1
2,4,5-Trichlorophenol	ND		2.1
2-Chloronaphthalene	ND		2.1
2-Nitroaniline	ND		11
Dimethyl phthalate	ND		5.3
Acenaphthylene	ND		2.1
3-Nitroaniline	ND		5.3
Acenaphthene	ND		2.1
2,4-Dinitrophenol	ND		11
4-Nitrophenol	ND		11
Dibenzofuran	ND		2.1
2,4-Dinitrotoluene	ND		2.1
2,6-Dinitrotoluene	ND		5.3
Diethyl phthalate	ND		5.3
4-Chlorophenyl phenyl ether	ND		5.3
Fluorene	ND		2.1
4-Nitroaniline	ND		11
2-Methyl-4,6-dinitrophenol	ND		11
N-Nitrosodiphenylamine	ND		2.1
· ·	· • 		—··

Client: LFR, Inc. Job Number: 720-18554-1

Client Sample ID: MW-3

Lab Sample ID: 720-18554-2 Date Sampled: 03/16/2009 1505 Client Matrix: Water 03/16/2009 1630 Date Received:

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

8270C Analysis Batch: 720-48005 Method: Instrument ID:

Preparation: 3510C Prep Batch: 720-47916 d:\data\200903\032509\720-Lab File ID:

Dilution: 1.0

Initial Weight/Volume: 950 mL Date Analyzed: 03/25/2009 1329 Final Weight/Volume: 1 mL Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result (ug/L)	Qualifier	RL
Hexachlorobenzene	ND		2.1
Pentachlorophenol	ND		11
Phenanthrene	ND		2.1
Anthracene	ND		2.1
Di-n-butyl phthalate	ND		5.3
Fluoranthene	ND		2.1
Pyrene	ND		2.1
Butyl benzyl phthalate	ND		5.3
3,3'-Dichlorobenzidine	ND		5.3
Benzo[a]anthracene	ND		5.3
Bis(2-ethylhexyl) phthalate	ND		11
Chrysene	ND		2.1
Di-n-octyl phthalate	ND		21
Benzo[b]fluoranthene	ND		2.1
Benzo[a]pyrene	ND		2.1
Benzo[k]fluoranthene	ND		2.1
Indeno[1,2,3-cd]pyrene	ND		2.1
Benzo[g,h,i]perylene	ND		2.1
Benzoic acid	ND		11
Azobenzene	ND		2.1
Dibenz(a,h)anthracene	ND		2.1
Surrogate	%Rec		Acceptance Limits
Nitrobenzene-d5	70		6 - 98
2-Fluorobiphenyl	69		6 - 103
Terphenyl-d14	78		36 - 106
2-Fluorophenol	39		1 - 66
Phenol-d5	28		1 - 47
2,4,6-Tribromophenol	80		22 - 124

DATA REPORTING QUALIFIERS

Lab Section Qualifier Description

Client: LFR, Inc. Job Number: 720-18554-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC/MS VOA	·				•
Analysis Batch:720-47847	7				
LCS 720-47847/2	Lab Control Sample	T	Water	8260B/CA_LUFT	
LCSD 720-47847/1	Lab Control Sample Duplicate	T	Water	8260B/CA_LUFT	
MB 720-47847/3	Method Blank	T	Water	8260B/CA_LUFT	
720-18554-1	MW-1	T	Water	8260B/CA_LUFT	
720-18554-2	MW-3	Т	Water	8260B/CA_LUFT	
Report Basis T = Total					
GC/MS Semi VOA					
Prep Batch: 720-47916					
LCS 720-47916/2-A	Lab Control Sample	T	Water	3510C	
LCSD 720-47916/3-A	Lab Control Sample Duplicate	T	Water	3510C	
MB 720-47916/1-A	Method Blank	T	Water	3510C	
720-18554-1	MW-1	T	Water	3510C	
720-18554-2	MW-3	Т	Water	3510C	
Analysis Batch:720-48005	5				
LCS 720-47916/2-A	Lab Control Sample	T	Water	8270C	720-47916
LCSD 720-47916/3-A	Lab Control Sample Duplicate	T	Water	8270C	720-47916
		Т	\A/a+a=	00700	700 47040
MB 720-47916/1-A	Method Blank	ı	Water	8270C	720-47916
MB 720-47916/1-A 720-18554-1	Method Blank MW-1	T	Water	8270C 8270C	720-47916 720-47916

Report Basis

T = Total

Client: LFR, Inc. Job Number: 720-18554-1

Method Blank - Batch: 720-47847 Method: 8260B/CA_LUFTMS

Preparation: 5030B

Lab Sample ID: MB 720-47847/3 Analysis Batch: 720-47847 Instrument ID: Varian 3900A

Client Matrix: Water Prep Batch: N/A Lab File ID: e:\data\2009\200903\03190

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL

Date Analyzed: 03/19/2009 1000 Final Weight/Volume: 10 mL Date Prepared: 03/19/2009 1000

Analyte	Result	Qual	RL
Benzene	ND		0.50
Gasoline Range Organics (GRO)-C5-C12	ND		50
TAME	ND		0.50
Ethyl tert-butyl ether	ND		0.50
Toluene	ND		0.50
Xylenes, Total	ND		1.0
MTBE	ND		0.50
EDB	ND		0.50
DIPE	ND		1.0
TBA	ND		5.0
1,2-DCA	ND		0.50
Ethylbenzene	ND		0.50
Surrogate	% Rec	Acceptance Limits	3
Toluene-d8 (Surr)	90	78 - 112	
1,2-Dichloroethane-d4 (Surr)	92	67 - 126	

Client: LFR, Inc. Job Number: 720-18554-1

Lab Control Sample/ Method: 8260B/CA_LUFTMS

Lab Control Sample Duplicate Recovery Report - Batch: 720-47847 Preparation: 5030B

LCS Lab Sample ID: LCS 720-47847/2 Analysis Batch: 720-47847 Instrument ID: Varian 3900A

Client Matrix: Water Prep Batch: N/A Lab File ID: e:\data\2009\200903\03190

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL Date Analyzed: 03/19/2009 1036 Final Weight/Volume: 10 mL

Date Prepared: 03/19/2009 1036

LCSD Lab Sample ID: LCSD 720-47847/1 Analysis Batch: 720-47847 Instrument ID: Varian 3900A

Date Prepared:

03/19/2009 1059

Client Matrix: Water Prep Batch: N/A Lab File ID: e:\data\2009\200903\031909

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 10 mL

Date Analyzed: 03/19/2009 1059 Final Weight/Volume: 10 mL

% Rec. LCS **LCSD RPD** RPD Limit LCS Qual LCSD Qual Analyte Limit 74 - 112 Benzene 84 89 5 20 Gasoline Range Organics (GRO)-C5-C12 63 42 - 80 20 61 4 Toluene 72 76 65 - 98 4 20 **MTBE** 84 69 - 104 4 20 80 Surrogate LCS % Rec LCSD % Rec Acceptance Limits 92 Toluene-d8 (Surr) 93 78 - 112 1,2-Dichloroethane-d4 (Surr) 90 84 67 - 126

Client: LFR, Inc. Job Number: 720-18554-1

Method Blank - Batch: 720-47916 Method: 8270C Preparation: 3510C

Lab Sample ID: MB 720-47916/1-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\mb

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL
Date Analyzed: 03/25/2009 1149

Date Prepared: 03/23/2009 1257

Units: ug/L Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume: 1.0 uL

Result Qual RL Analyte Phenol ND 2.0 Bis(2-chloroethyl)ether ND 2.0 2-Chlorophenol ND 2.0 1,3-Dichlorobenzene ND 2.0 1,4-Dichlorobenzene ND 2.0 Benzyl alcohol 5.0 ND 1,2-Dichlorobenzene ND 2.0 2-Methylphenol ND 2.0 4-Methylphenol 2.0 ND N-Nitrosodi-n-propylamine ND 2.0 Hexachloroethane ND 2.0 Nitrobenzene 2.0 ND 2.0 Isophorone ND 2-Nitrophenol ND 2.0 2.0 2,4-Dimethylphenol ND Bis(2-chloroethoxy)methane 5.0 ND 5.0 2,4-Dichlorophenol ND 1,2,4-Trichlorobenzene 2.0 ND Naphthalene ND 2.0 4-Chloroaniline ND 2.0 Hexachlorobutadiene ND 2.0 4-Chloro-3-methylphenol ND 5.0 2-Methylnaphthalene ND 2.0 Hexachlorocyclopentadiene ND 5.0 2,4,6-Trichlorophenol ND 2.0 2,4,5-Trichlorophenol ND 2.0 2-Chloronaphthalene ND 2.0 2-Nitroaniline ND 10 Dimethyl phthalate ND 5.0 Acenaphthylene ND 2.0 3-Nitroaniline ND 5.0 Acenaphthene ND 2.0 2,4-Dinitrophenol ND 10 4-Nitrophenol ND 10 Dibenzofuran ND 2.0 2.4-Dinitrotoluene 2.0 ND 2,6-Dinitrotoluene ND 5.0 Diethyl phthalate ND 5.0 4-Chlorophenyl phenyl ether 5.0 ND Fluorene ND 2.0 4-Nitroaniline ND 10

Client: LFR, Inc. Job Number: 720-18554-1

Method Blank - Batch: 720-47916 Method: 8270C Preparation: 3510C

Lab Sample ID: MB 720-47916/1-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\mb

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL Date Analyzed: 03/25/2009 1149 Final Weight/Volume: 1 mL

Date Analyzed: 03/25/2009 1149 Final Weight/Volume: 1 mL

Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

Analyte	Result	Qual	RL
2-Methyl-4,6-dinitrophenol	ND		10
N-Nitrosodiphenylamine	ND		2.0
4-Bromophenyl phenyl ether	ND		5.0
Hexachlorobenzene	ND		2.0
Pentachlorophenol	ND		10
Phenanthrene	ND		2.0
Anthracene	ND		2.0
Di-n-butyl phthalate	ND		5.0
Fluoranthene	ND		2.0
Pyrene	ND		2.0
Butyl benzyl phthalate	ND		5.0
3,3'-Dichlorobenzidine	ND		5.0
Benzo[a]anthracene	ND		5.0
Bis(2-ethylhexyl) phthalate	ND		10
Chrysene	ND		2.0
Di-n-octyl phthalate	ND		20
Benzo[b]fluoranthene	ND		2.0
Benzo[a]pyrene	ND		2.0
Benzo[k]fluoranthene	ND		2.0
Indeno[1,2,3-cd]pyrene	ND		2.0
Benzo[g,h,i]perylene	ND		2.0
Benzoic acid	ND		10
Azobenzene	ND		2.0
Dibenz(a,h)anthracene	ND		2.0
Surrogate	% Rec	Acceptance Limits	
Nitrobenzene-d5	72	6 - 98	
2-Fluorobiphenyl	69	6 - 103	
Terphenyl-d14	84	36 - 106	
2-Fluorophenol	46	1 - 66	
Phenol-d5	32	1 - 47	
2,4,6-Tribromophenol	74	22 - 124	

Client: LFR, Inc. Job Number: 720-18554-1

Lab Control Sample/ Method: 8270C Lab Control Sample Duplicate Recovery Report - Batch: 720-47916 Preparation: 3510C

LCS Lab Sample ID: LCS 720-47916/2-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Prep Batch: 720-47916 d:\data\200903\032509\LC Water Lab File ID:

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL Date Analyzed: 03/25/2009 1042 Final Weight/Volume: 1 mL

Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

LCSD Lab Sample ID: LCSD 720-47916/3-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\LCS

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL Date Analyzed: 03/25/2009 1116 Final Weight/Volume: 1 mL Injection Volume: Date Prepared: 03/23/2009 1257 1.0 uL

	<u>%</u>	Rec.					
Analyte	LCS	LCSD	Limit	RPD	RPD Limit	LCS Qual	LCSD Qual
Phenol	31	38	10 - 49	22	35		
Bis(2-chloroethyl)ether	58	71	22 - 113	21	35		
2-Chlorophenol	60	76	19 - 104	24	25		
1,3-Dichlorobenzene	62	70	18 - 95	13	35		
1,4-Dichlorobenzene	54	67	17 - 82	22	30		
Benzyl alcohol	60	75	21 - 94	22	35		
1,2-Dichlorobenzene	64	71	20 - 89	10	35		
2-Methylphenol	58	70	18 - 100	19	35		
4-Methylphenol	51	62	16 - 93	19	35		
N-Nitrosodi-n-propylamine	64	81	34 - 108	23	34		
Hexachloroethane	64	75	4 - 87	16	35		
Nitrobenzene	74	96	27 - 117	26	35		
Isophorone	79	95	42 - 118	19	35		
2-Nitrophenol	67	82	34 - 107	20	35		
2,4-Dimethylphenol	81	98	42 - 122	19	35		
Bis(2-chloroethoxy)methane	65	77	37 - 110	18	35		
2,4-Dichlorophenol	67	81	39 - 107	20	35		
1,2,4-Trichlorobenzene	62	74	20 - 95	17	35		
Naphthalene	65	80	22 - 99	21	35		
4-Chloroaniline	45	54	10 - 80	17	35		
Hexachlorobutadiene	52	66	3 - 101	24	35		
4-Chloro-3-methylphenol	74	88	22 - 147	17	31		
2-Methylnaphthalene	64	77	10 - 130	19	35		
Hexachlorocyclopentadiene	65	79	21 - 101	19	35		
2,4,6-Trichlorophenol	72	81	31 - 118	12	35		
2,4,5-Trichlorophenol	71	83	35 - 112	16	35		
2-Chloronaphthalene	63	79	26 - 106	22	35		
2-Nitroaniline	72	86	31 - 106	18	35		
Dimethyl phthalate	72	88	47 - 117	20	35		
Acenaphthylene	83	103	33 - 128	22	35		
3-Nitroaniline	71	85	40 - 107	18	35		
Acenaphthene	75	81	31 - 109	8	30		

Client: LFR, Inc. Job Number: 720-18554-1

Lab Control Sample/
Lab Control Sample Duplicate Recovery Report - Batch: 720-47916

Method: 8270C
Preparation: 3510C

LCS Lab Sample ID: LCS 720-47916/2-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\LC

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL Date Analyzed: 03/25/2009 1042 Final Weight/Volume: 1 mL

Date Prepared: 03/23/2009 1257 Injection Volume: 1.0 uL

LCSD Lab Sample ID: LCSD 720-47916/3-A Analysis Batch: 720-48005 Instrument ID: Sat 2K1

Client Matrix: Water Prep Batch: 720-47916 Lab File ID: d:\data\200903\032509\LCS

Dilution: 1.0 Units: ug/L Initial Weight/Volume: 1000 mL

 Date Analyzed:
 03/25/2009 1116

 Date Prepared:
 03/23/2009 1257

 Injection Volume:
 1.0 uL

	0	% Rec.					
Analyte	LCS	LCSD	Limit	RPD	RPD Limit	LCS Qual	LCSD Qual
2,4-Dinitrophenol	67	76	51 - 102	13	35		
4-Nitrophenol	42	48	26 - 63	13	35		
Dibenzofuran	68	79	42 - 97	14	35		
2,4-Dinitrotoluene	77	92	52 - 113	18	35		
2,6-Dinitrotoluene	77	83	50 - 115	8	35		
Diethyl phthalate	76	89	54 - 111	16	35		
4-Chlorophenyl phenyl ether	73	84	40 - 110	14	35		
Fluorene	71	85	37 - 104	17	35		
4-Nitroaniline	79	89	56 - 116	12	35		
2-Methyl-4,6-dinitrophenol	66	82	47 - 114	21	35		
N-Nitrosodiphenylamine	80	96	56 - 123	19	35		
4-Bromophenyl phenyl ether	76	95	52 - 111	22	35		
Hexachlorobenzene	78	93	61 - 104	18	35		
Pentachlorophenol	67	80	55 - 107	18	35		
Phenanthrene	75	91	56 - 110	19	35		
Anthracene	80	94	58 - 114	16	35		
Di-n-butyl phthalate	86	104	56 - 114	19	35		
Fluoranthene	83	96	60 - 121	15	35		
Pyrene	76	83	56 - 91	10	35		
Butyl benzyl phthalate	73	82	37 - 100	12	35		
3,3'-Dichlorobenzidine	59	68	37 - 111	15	35		
Benzo[a]anthracene	67	79	50 - 112	16	35		
Bis(2-ethylhexyl) phthalate	77	88	59 - 111	13	35		
Chrysene	73	77	56 - 94	4	35		
Di-n-octyl phthalate	71	83	47 - 118	15	35		
Benzo[b]fluoranthene	75	80	55 - 110	7	35		
Benzo[a]pyrene	66	75	59 - 103	12	35		
Benzo[k]fluoranthene	73	89	55 - 110	20	35		
Indeno[1,2,3-cd]pyrene	78	88	63 - 126	11	35		
Benzo[g,h,i]perylene	79	89	10 - 140	12	35		
Benzoic acid	28	30	7 - 46	8	35		
Azobenzene	76	87	49 - 102	13	35		

Client: LFR, Inc. Job Number: 720-18554-1

Lab Control Sample/ Method: 8270C Lab Control Sample Duplicate Recovery Report - Batch: 720-47916 Preparation: 3510C

LCS Lab Sample ID: LCS 720-47916/2-A

Client Matrix: Water

Dilution: 1.0

Date Analyzed: 03/25/2009 1042 Date Prepared: 03/23/2009 1257 Analysis Batch: 720-48005 Prep Batch: 720-47916

Units: ug/L

Instrument ID: Sat 2K1

d:\data\200903\032509\LC Lab File ID:

Initial Weight/Volume: 1000 mL Final Weight/Volume: 1 mL Injection Volume: 1.0 uL

LCSD Lab Sample ID: LCSD 720-47916/3-A

Client Matrix: Water Dilution: 1.0

Date Analyzed: 03/25/2009 1116 Date Prepared: 03/23/2009 1257

Analysis Batch: 720-48005

Prep Batch: 720-47916

Units: ug/L

Instrument ID: Sat 2K1

Lab File ID: d:\data\200903\032509\LCS

Initial Weight/Volume: 1000 mL Final Weight/Volume: 1 mL Injection Volume: 1.0 uL

	9	<u>% Rec.</u>					
Analyte	LCS	LCSD	Limit	RPD	RPD Limit	LCS Qual	LCSD Qual
Dibenz(a,h)anthracene	69	82	10 - 130	17	35		
Surrogate	L	CS % Rec	LCSD %	Rec	Accep	tance Limits	:
Nitrobenzene-d5	7	70	86		6	- 98	
2-Fluorobiphenyl	6	64	76		6	- 103	
Terphenyl-d14	7	74	83		3	6 - 106	
2-Fluorophenol	4	10	47		1	- 66	
Phenol-d5	2	29	35		1	- 47	
2,4,6-Tribromophenol	7	75	83		2	2 - 124	

SAMPLE COLLECTOR:	Street, 12th	n Floor	PROJEC	04567	-07	SECTION	ON NO.:		DATE:	3/19	69	SAMP	PZ CINITIALS:	SERIAL NO	
Emeryville, C	alifornia 94	608- 10) 652-2246	PROJEC	t NAME: fahSat		adia	M		SAMPLE	R (Sign	ature):	Tone	w-	Nº 2	03315
			/PLE		100			16		А	NALYS	SESn	15		REMARKS
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MW-3	3/16	15:05	6	X	X	×	$\times \times$		×	×	X	X			
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Login Sample Receipt Check List

Client: LFR, Inc. Job Number: 720-18554-1

Login Number: 18554 List Source: TestAmerica San Francisco

Creator: Bullock, Tracy List Number: 1

Question	T / F/ NA	Comment
Radioactivity either was not measured or, if measured, is at or below background	N/A	
The cooler's custody seal, if present, is intact.	N/A	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the sample IDs on the containers and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	False	see NCM
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	



ANALYTICAL REPORT

Job Number: 720-18554-3

Job Description: Hanson Radum

For: LFR, Inc. 1900 Powell St 12th Floor Emeryville, CA 94608-1827

Attention: Mr. Ron Goloubow

Approved for release Afsaneh Salimpour Project Manager I 3/31/2009 3:17 PM

Afsaneh Salimpour Project Manager I afsaneh.salimpour@testamericainc.com 03/31/2009

Akanef Sal

EXECUTIVE SUMMARY - Detections

Client: LFR, Inc. Job Number: 720-18554-3

Lab Sample ID Client Sample ID Reporting
Analyte Result / Qualifier Limit Units Method

No Detections

METHOD SUMMARY

Client: LFR, Inc. Job Number: 720-18554-3

Description	Lab Location	Method	Preparation Method
Matrix: Water			
Diesel Range Organics (DRO) (GC)	TAL SF	SW846 8015B	
Liquid-Liquid Extraction (Separatory Funnel)	TAL SF		SW846 3510C SGC

Lab References:

TAL SF = TestAmerica San Francisco

Method References:

SW846 = "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

METHOD / ANALYST SUMMARY

Client: LFR, Inc. Job Number: 720-18554-3

 Method
 Analyst
 Analyst ID

 SW846
 8015B
 Hayashi, Derek
 DH

SAMPLE SUMMARY

Client: LFR, Inc. Job Number: 720-18554-3

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received	
720-18554-1	MW-1	Water	03/16/2009 1440	03/16/2009 1630	
720-18554-2	MW-3	Water	03/16/2009 1505	03/16/2009 1630	

Analytical Data

Client: LFR, Inc. Job Number: 720-18554-3

Client Sample ID: MW-1

Lab Sample ID: 720-18554-1 Date Sampled: 03/16/2009 1440 Client Matrix: Water Date Received: 03/16/2009 1630

8015B Diesel Range Organics (DRO) (GC)-Silica Gel Cleanup

8015B Analysis Batch: 720-48168 Instrument ID: HP DRO5 Method: Preparation:

3510C SGC Prep Batch: 720-48032 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume: 500 mL Date Analyzed: 03/27/2009 0140 Final Weight/Volume: 2 mL

Date Prepared: 03/26/2009 1221 Injection Volume:

Column ID: **PRIMARY**

Analyte	Result (ug/L)	Qualifier	RL
Motor Oil Range Organics [C24-C36]	ND		300
Diesel Range Organics [C9-C24]	ND		50

%Rec Acceptance Limits Surrogate 0 - 5 Capric Acid (Surr) 0 p-Terphenyl 78 31 - 120

Analytical Data

Client: LFR, Inc. Job Number: 720-18554-3

Client Sample ID: MW-3

Lab Sample ID: 720-18554-2 Date Sampled: 03/16/2009 1505 Client Matrix: Water Date Received: 03/16/2009 1630

8015B Diesel Range Organics (DRO) (GC)-Silica Gel Cleanup

Analysis Batch: 720-48168 HP DRO5 Method: 8015B Instrument ID: Preparation: 3510C SGC Prep Batch: 720-48032 Lab File ID: N/A

Dilution: 1.0 Initial Weight/Volume:

500 mL Date Analyzed: 03/27/2009 0207 Final Weight/Volume: 2 mL

Date Prepared: 03/26/2009 1221 Injection Volume:

Column ID: **PRIMARY**

Analyte Result (ug/L) Qualifier RLMotor Oil Range Organics [C24-C36] ND 300 Diesel Range Organics [C9-C24] ND 50

Surrogate %Rec Acceptance Limits 0 - 5 Capric Acid (Surr) 0 p-Terphenyl 85 31 - 120

DATA REPORTING QUALIFIERS

Lab Section Qualifier Description

Client: LFR, Inc. Job Number: 720-18554-3

QC Association Summary

		Report			
Lab Sample ID	Client Sample ID	Basis	Client Matrix	Method	Prep Batch
GC Semi VOA					
Prep Batch: 720-48032					
LCS 720-48032/2-A	Lab Control Spike	Α	Water	3510C SGC	
_CSD 720-48032/3-A	Lab Control Spike Duplicate	Α	Water	3510C SGC	
MB 720-48032/1-A	Method Blank	Α	Water	3510C SGC	
720-18554-1	MW-1	Α	Water	3510C SGC	
720-18554-2	MW-3	Α	Water	3510C SGC	
Analysis Batch:720-48	168				
LCS 720-48032/2-A	Lab Control Spike	Α	Water	8015B	720-48032
_CSD 720-48032/3-A	Lab Control Spike Duplicate	Α	Water	8015B	720-48032
MB 720-48032/1-A	Method Blank	Α	Water	8015B	720-48032
720-18554-1	MW-1	Α	Water	8015B	720-48032
720-18554-2	MW-3	Α	Water	8015B	720-48032

Report Basis

A = Silica Gel Cleanup

Client: LFR, Inc. Job Number: 720-18554-3

Method Blank - Batch: 720-48032 Method: 8015B

Preparation: 3510C SGC Silica Gel Cleanup

Lab Sample ID: MB 720-48032/1-A

Client Matrix: Water Dilution: 1.0

Date Analyzed: 03/27/2009 0112 Date Prepared: 03/26/2009 1221 Analysis Batch: 720-48168 Prep Batch: 720-48032

Units: ug/L

Instrument ID: HP DRO5 Lab File ID: N/A

Initial Weight/Volume: 500 mL Final Weight/Volume: 2 mL

Injection Volume:

Column ID: PRIMARY

Analyte	Result	Qual	RL
Motor Oil Range Organics [C24-C36] Diesel Range Organics [C9-C24]	ND ND		300 50
Surrogate	% Rec	Acceptance Limits	
Capric Acid (Surr) p-Terphenyl	0 85	0 - 5 31 - 120	

Lab Control Spike/ Method: 8015B

Lab Control Spike Duplicate Recovery Report - Batch: 720-48032 Preparation: 3510C SGC

Silica Gel Cleanup

LCS Lab Sample ID: LCS 720-48032/2-A Analysis Batch: 720-48168 Instrument ID: HP DRO5

Client Matrix: Water Prep Batch: 720-48032 Lab File ID: N/A
Dilution: 1.0 Units: ug/L Initial Weight/Volume:

Date Analyzed: 03/27/2009 0017 Final Weight/Volume: 2 mL
Date Prepared: 03/26/2009 1221 Injection Volume:
Column ID: PRIMARY

LCSD Lab Sample ID: LCSD 720-48032/3-A

Client Matrix: Water Dilution: 1.0

Date Analyzed: 03/27/2009 0045 Date Prepared: 03/26/2009 1221 Analysis Batch: 720-48168 Prep Batch: 720-48032

Inite: us/l

Units: ug/L

Instrument ID: HP DRO5

Lab File ID: N/A

Initial Weight/Volume: 500 mL Final Weight/Volume: 2 mL

500 mL

Injection Volume:

Column ID: PRIMARY

	<u>9</u>	<u> 6 Rec.</u>					
Analyte	LCS	LCSD	Limit	RPD	RPD Limit	LCS Qual	LCSD Qual
Diesel Range Organics [C9-C24]	78	105	49 - 120	29	30		
Surrogate	L	CS % Rec	LCSD %	Rec	Accep	tance Limits	;
p-Terphenyl	9	5	93		3	1 - 120	

AMPLE COLLECTOR:	Street 12th		PROJEC	04567			ON NO.:		DATE				AMPLET 7	Z NITIALS:	SERIAL	//500 No.:
Emeryville, C (510) 652-45	California 94	608	PROJEC	TNAME:		adu	M		SAMP	LER (Si	gnature)	100	-6		Nº	20331
			IPLE		- 17			/a			ANAL		1	15		REMARK:
120-185	=355.00	4	Lab Sample N	of Containers		TYPE	Prino Esas	BIET UC	BOZ 18OZ	Soleta IV	Fur	0+4	cayer	nato vi o	B260	OCs: **Metals List CAM List RCR
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MW-3	3/16	15:05	6	X	X	X	XX		- 1	XX	- X		X			_
161-3-4	3/11	15:40	4	X	X	×	X					-	X			
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Login Sample Receipt Check List

Client: LFR, Inc. Job Number: 720-18554-3

Login Number: 18554 List Source: TestAmerica San Francisco

Creator: Bullock, Tracy List Number: 1

Question	T / F/ NA	Comment
Radioactivity either was not measured or, if measured, is at or below background	N/A	
The cooler's custody seal, if present, is intact.	N/A	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the sample IDs on the containers and the COC.	True	
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	False	see NCM
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	