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November 7, 2014

Alameda County Department of
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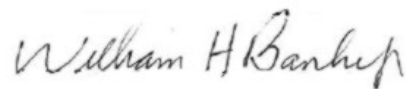
Attention: Mark Detterman

Subject: Third Quarter 2014 Groundwater Monitoring Report
3800 San Pablo Avenue, Emeryville, California
ACDEH Fuel Leak Case: RO00002520; Global ID: T06019788682

Ladies and Gentlemen:

Attached please find a copy of the *Third Quarter 2014 Groundwater Monitoring* prepared by Gribi Associates. I declare, under penalty of perjury, that the information and/or recommendations contained in the attached document or report is true and correct to the best of my knowledge.

Very truly yours,



William H. Banker, Jr.
San Pablo Avenue Venture
c/o Banker, Marks & Kirk
1720 Broadway, Suite 202
Oakland, CA 94612



November 7, 2014

Alameda County Department of
Environmental Health
1131 Harbor Bay Parkway, 2nd Floor
Alameda, CA 94502

Attention: Mark Detterman

Subject: Third Quarter 2014 Groundwater Monitoring Report
3800 San Pablo Avenue, Emeryville, California
ACDEH Fuel Leak Case: RO00002520; Global ID: T06019788682

Ladies and Gentlemen:

Gribi Associates is pleased to submit this *Third Quarter 2014 Groundwater Monitoring Report* on behalf San Pablo Avenue Venture for the property located at 3800 San Pablo Avenue in Emeryville, California (see Figure 1 and Figure 2). This letter report documents the monitoring and sampling of four site wells on September 29, 2014.

DESCRIPTION OF SAMPLING ACTIVITIES

1. Gribi Associates personnel conducted groundwater monitoring and sampling activities for four site wells (MW-1, MW-2, MW-3, and MW-4) on September 29, 2014.
2. Groundwater monitoring and sampling was conducted in accordance with California LUFT Field Manual, including the following:
 - a. measuring static water levels;
 - b. checking for presence of free-product;
 - c. and purging of approximately three well volumes while recording of temperature, pH, conductivity, and clarity.
3. Collected groundwater samples were placed in an ice-chilled cooler and submitted to a state-certified laboratory for analyses.
4. Copies of groundwater sampling field data sheets are provided as Attachment A.

RESULTS OF GROUNDWATER MONITORING

Hydrologic Conditions

1. Groundwater depths ranged from approximately 10.31 feet (MW-3) to 11.28 feet (MW-2).
2. Groundwater elevations ranged from 27.29 feet above means sea level (msl) (MW-4) to 28.53 feet msl (MW-3).
3. Groundwater potentiometric gradient during this monitoring event was to the east at an approximate gradient of 0.1 feet/feet.
4. Groundwater elevations and contours are shown on Figure 3.

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2. Groundwater elevations ranged from 27.29 feet above means sea level (msl) (MW-4) to 28.53 feet msl (MW-3).
3. Groundwater potentiometric gradient during this monitoring event was to the east at an approximate gradient of 0.1 feet/foot.
4. Groundwater elevations and contours are shown on Figure 3.

Laboratory Analytical Results

1. Groundwater samples from the four sampled wells were analyzed for the following parameters with standard method turn-around-time on results:
 - a. USEPA 8260B Total Petroleum Hydrocarbons as Gasoline (TPH-G)
 - b. USEPA 8260B Benzene, Toluene, Ethylbenzene, and Xylenes (BTEX)
 - c. USEPA 8260B Oxygenates (DIPE, ETBE, MTBE, TAME, TBA)
 - d. USEPA 8260B Lead Scavengers
 - e. USEPA 8260B Volatile Organic Compounds (VOCs)
 - f. USEPA 8260B Naphthalene
 - g. USEPA 8015C Total Petroleum Hydrocarbons as Diesel (TPH-D)
 - h. USEPA 8015C Total Petroleum Hydrocarbons as Heating Oil (TPH-HO)
 - i. USEPA 8270C Semi-Volatile Organic Compounds (SVOCs)
2. Groundwater analytical results are summarized in Table 1 and on Figure 4.
3. Groundwater hydrocarbon trends for selected wells are provided as Attachment B.
4. The laboratory analytical data report and chain-of custody are provided as Attachment C.

SITE REMEDIATION ACTIVITIES

1. Gribi Associates installed an ozone remediation system at the site during the week of September 2, 2013.
2. The ozone system was started on September 9, 2013.
 - a. The system operated continuously until the mid-October 2013.
 - b. The system required repairs and was re-started on November 7, 2013 and operated continuously until the system was turned off on January 17, 2014.
3. Gribi Associates resumed ozone remediation at the site on August 5, 2014 and turned it off on October 24, 2014.

CONCLUSIONS

1. Following re-implementation of the ozone remediation system, groundwater hydrocarbon concentrations dropped to relatively low levels in wells MW-1, MW-2 and MW-3.
 - a. At MW-1, respective groundwater TPH-G and Benzene concentrations decreased from 8,600 ug/L and 880 ug/L in June 2013 to 400 ug/L and nondetect during this monitoring event.

- b. At MW-2, respective groundwater TPH-G and Benzene concentrations decreased from 12,000 ug/L and 870 ug/L in June 2013 to nondetect and 2.3 ug/L during this monitoring event.
 - c. At MW-3, groundwater TPH-G and Benzene concentrations decreased from 12,000 ug/L and 1,400 ug/L in June 2013 to 400 ug/L and nondetect during this monitoring event.
 - d. At MW-4, groundwater TPH-G and Benzene concentrations decreased from 6,300 ug/L and 72 ug/L in June 2013 to 5,600 ug/L and 16 ug/L during this monitoring event.
2. Groundwater samples from wells MW-1 and MW-4 showed relatively low concentrations of Diesel and Heating Oil range hydrocarbons.

PLANNED ACTIVITIES

1. Gribi Associates plans to conduct a quarterly groundwater and soil vapor monitoring and sampling event during the fourth quarter of 2014.

We appreciate this opportunity to provide this report for your review. Please contact us if there are questions or if additional information is required.

Very truly yours,



Matthew A. Rosman
Project Engineer



James E. Gribi
Professional Geologist
California No. 5843



Enclosure

- c: Mr. Bill Banker, Jr., San Pablo Avenue Venture

TABLE

Table 1
CUMULATIVE GROUNDWATER LABORATORY ANALYTICAL RESULTS
Former Maz Glass UST Site

Well ID	Date	GW Depth	GW Elev.	Groundwater Concentration, in micrograms per liter (ug/L)												
				TPH-G	TPH-D	TPH-HO	B	T	E	X	OXY	Cr6	Br	N	SVOCs	Other VOCs
MW-1	5/18/2012	8.42	30.54	17,000	-	-	1,300	29	770	260	All ND	-	-	-	-	-
<38.96>	9/13/2012	10.55	28.41	13,000	-	-	630	10	780	86.7	All ND	-	-	-	-	-
	11/9/2012	9.72	29.24	15,000	-	-	1,200	21	1,100	283	All ND	-	-	-	-	-
	2/20/2013	8.34	30.62	9,800	-	-	970	15	860	171.5	All ND	-	-	75	-	-
	6/4/2013	9.39	29.57	8,600	-	-	880	15	770	121.2	All ND	-	-	74	-	-
Ozone Injection Started on September 9, 2013																
	9/26/2013	10.38	28.58	16,000	-	-	220	8.9	610	152.4	All ND	<0.20	0.091	120	-	-
	12/30/2013	9.92	29.04	4,700	-	-	62	1.5	110	62.75	All ND	-	-	23	-	-
Ozone Injection Stopped on February 7, 2014																
	3/7/2014	6.56	32.40	5,600	-	-	320	8.4	370	89.7	All ND	<0.20	0.047	68	-	-
	5/27/2014	9.77	29.19	2,900	-	-	180	4.3	290	38.51	All ND	-	-	24	-	-
Ozone Injection Resumed on August 5, 2014																
	9/29/2014	11.25	27.71	400	<500	960	<0.50	<0.50	1.1	1.3	38 TBA	-	-	<1.0	ALL ND	7.0 1,3,5-Trimethylbenzene 4.3 1,2,4-Trimethylbenzene
MW-2	5/18/2012	8.78	30.18	10,000	-	-	610	26	340	69	All ND	-	-	-	-	-
<38.96>	9/13/2012	10.64	28.32	11,000	-	-	990	27	460	42.9	All ND	-	-	-	-	-
	11/9/2012	9.57	29.39	17,000	-	-	750	19	280	64.9	All ND	-	-	-	-	-
	2/20/2013	8.86	30.1	8,200	-	-	860	29	410	70	All ND	-	-	29	-	-
	6/4/2013	9.86	29.1	12,000	-	-	870	23	410	43.8	All ND	-	-	46	-	-
Ozone Injection Started on September 9, 2013																
	9/26/2013	13.32	25.64	930	-	-	39	5.6	26	20	All ND	1.1	0.09	13	-	-
	12/30/2013	10.33	28.63	270	-	-	7.9	<0.50	2.9	<1.0	TBA=20	-	-	<1.0	-	-
Ozone Injection Stopped on February 7, 2014																
	3/7/2014	6.95	32.01	440	-	-	41	0.91	4.2	2.9	All ND	<0.20	0.13	4.2	-	-
	5/27/2014	9.95	29.01	1,200	-	-	250	5.9	34	14.2	All ND	-	-	8.1	-	-
Ozone Injection Resumed on August 5, 2014																
	9/29/2014	11.28	27.68	180	<500	<500	4.5	<0.50	0.73	<1.0	87 TBA	-	-	<1.0	ALL ND	ALL ND
MW-3	5/18/2012	8.61	30.23	13,000	-	-	1,400	36	350	378	All ND	-	-	-	-	-
<38.84>	9/13/2012	10.3	28.54	12,000	-	-	1,800	25	680	565.5	All ND	-	-	-	-	-
	11/9/2012	9.25	29.59	17,000	-	-	2,000	32	540	318.6	All ND	-	-	-	-	-
	2/20/2013	8.8	30.04	12,000	-	-	1,400	15	330	43.9	All ND	-	-	8.4	-	-
	6/4/2013	9.49	29.35	12,000	-	-	1,400	11	89	32.4	All ND	-	-	13	-	-
Ozone Injection Started on September 9, 2013																
	9/26/2013	10.89	27.95	5,500	-	-	190	2.8	42	27	All ND	<0.20	0.096	18	-	-
	12/30/2013	14.59	24.25	380	-	-	8.3	<0.50	2.3	1.6	All ND	-	-	<1.0	-	-

Table 1 CUMULATIVE GROUNDWATER LABORATORY ANALYTICAL RESULTS Former Maz Glass UST Site																
Well ID	Date	GW Depth	GW Elev.	Groundwater Concentration, in micrograms per liter (ug/L)												
				TPH-G	TPH-D	TPH-HO	B	T	E	X	OXY	Cr6	Br	N	SVOCs	Other VOCs
Ozone Injection Stopped on February 7, 2014																
	3/7/2014	6.99	31.85	400	-	-	31	0.75	2.6	2.9	All ND	<0.20	0.083	1.9	-	-
	5/27/2014	9.63	29.21	510	-	-	120	1.3	9.8	2.8	All ND	-	-	<1.0	-	-
Ozone Injection Resumed on August 5, 2014																
	9/29/2014	10.31	28.53	<50	<500	<500	2.3	<0.50	<0.50	<1.0	All ND	-	-	<1.0	ALL ND	ALL ND
MW-4	5/18/2012	8.28	30.2	10,000	-	-	82	32	330	278	All ND	-	-	-	-	-
<38.48>	9/13/2012	8.8	29.68	10,000	-	-	110	24	270	178.1	All ND	-	-	-	-	-
	11/9/2012	8.06	30.42	11,000	-	-	110	13	170	124.4	All ND	-	-	-	-	-
	2/20/2013	8.16	30.32	4,500	-	-	100	9.5	190	65.3	All ND	-	-	7.1	-	-
	6/4/2013	8.73	29.75	6,300	-	-	72	6.2	61	48.4	All ND	-	-	12	-	-
Ozone Injection Started on September 9, 2013																
	9/26/2013	9.76	28.72	12,000	-	-	48	3.7	70	18.2	All ND	<0.20	0.056	13	-	-
	12/30/2013	9.81	28.67	7,600	-	-	50	6.6	68	104.3	All ND	-	-	37	-	-
Ozone Injection Stopped on February 7, 2014																
	3/7/2014	6.76	31.72	3,100	-	-	38	4.3	51	76.5	All ND	<0.020	0.016	20	-	-
	5/27/2014	9.11	29.37	2,900	-	-	47	3.5	68	68.6	All ND	-	-	<1.0	-	-
Ozone Injection Resumed on August 5, 2014																
	9/29/2014	11.19	27.29	5,600	2,200	4,900	16	0.78	6.1	9.04	All ND	-	-	<1.0	All ND	1.3 sec-Butylbenzene 2.8 Isopropylbenzene 2.9 p-Isopropylbenzene 5.7 n-Propylbenzene 22 1,3,5-Trimethylbenzene 20 1,2,4-Trimethylbenzene
Environmental Screening Levels				100	110	NL	27	95,000	310	37,000	110 TBA	21	NL	160	Various	Various

TABLE NOTES

GW Elev = Groundwater mean sea level elevation
 TPH-G = Total Petroleum Hydrocarbons as gasoline

B = Benzene,
 T = Toluene
 E = Ethylbenzene

TPH-D
 TPH-K

X = Xylenes
 OXY = Oxygenates, including MTBE = Methyl-t-Butyl Ether, ter-Butanol (TBA), Di-isopropyl Ether (DIPE), Ethyl-t-butyl Ether (ETBE), and Tert-amyl Methyl Ether (TAME).

Cr6 = Hexavalent Chromium

Br = Bromate
 N = Naphthalene.

<38.96> = Top of casing mean sea level elevation (Virgil Chavez Land Survey).

All ND = No detectable concentrations of all analytes.

- = Not analyzed for this analyte.

SVOCs = semi-volatile organic compounds

VOCs = volatile organic compounds

<1.0 = Not detected above the expressed value.

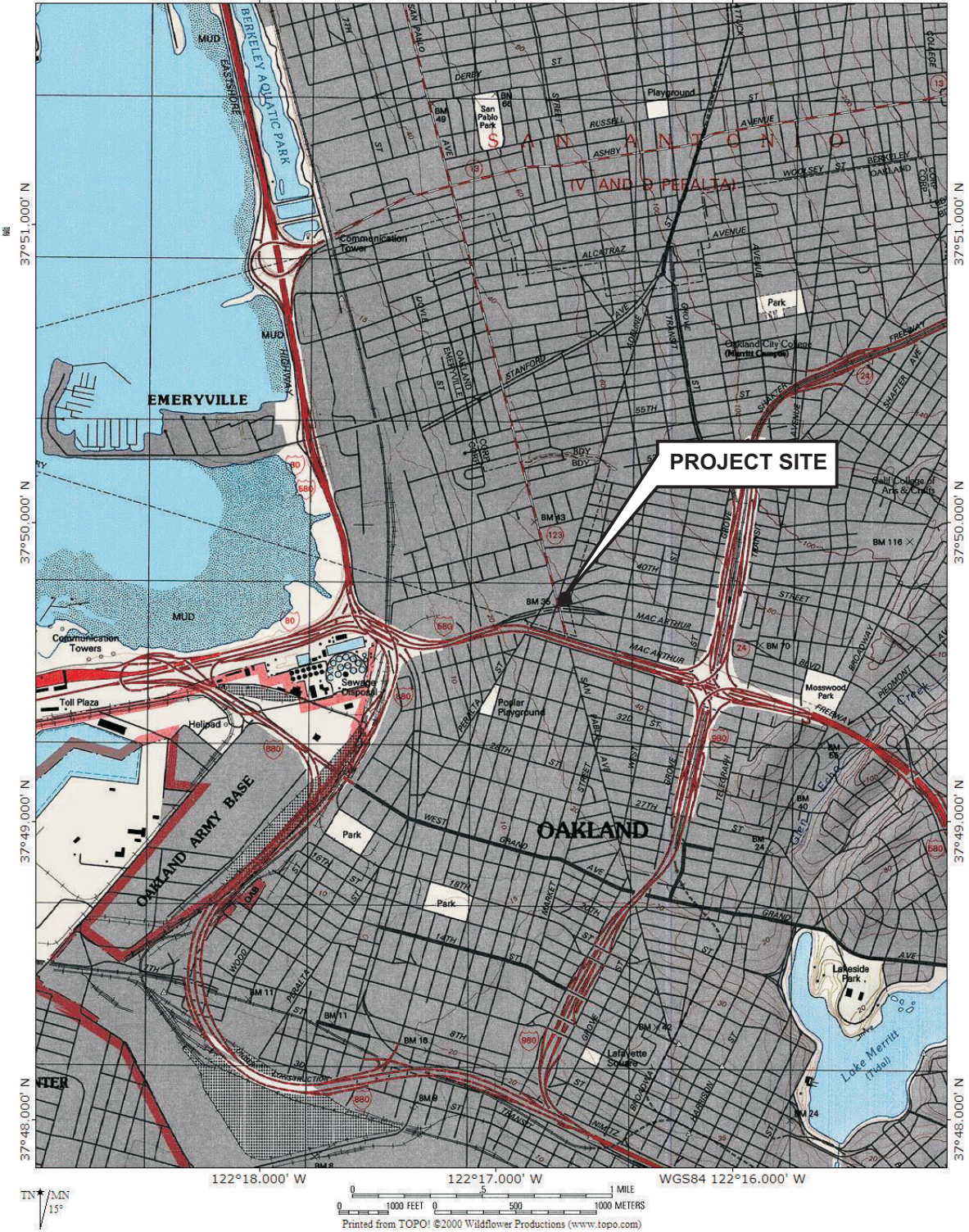
ESL = Environmental Screening Levels, as contained in *Screening for Environmental*

Concerns at Sites with Contaminated Soil and Groundwater, San Francisco Bay Regional Water Quality Control Board, December 2013, Table E-1, Groundwater to Indoor Air, fine grained soils, residential land use.

NL = Not Listed

FIGURES

TOPO! map printed on 04/03/07 from "California.tpo" and "Untitled.tpg"
 122°18.000' W 122°17.000' W WGS84 122°16.000' W



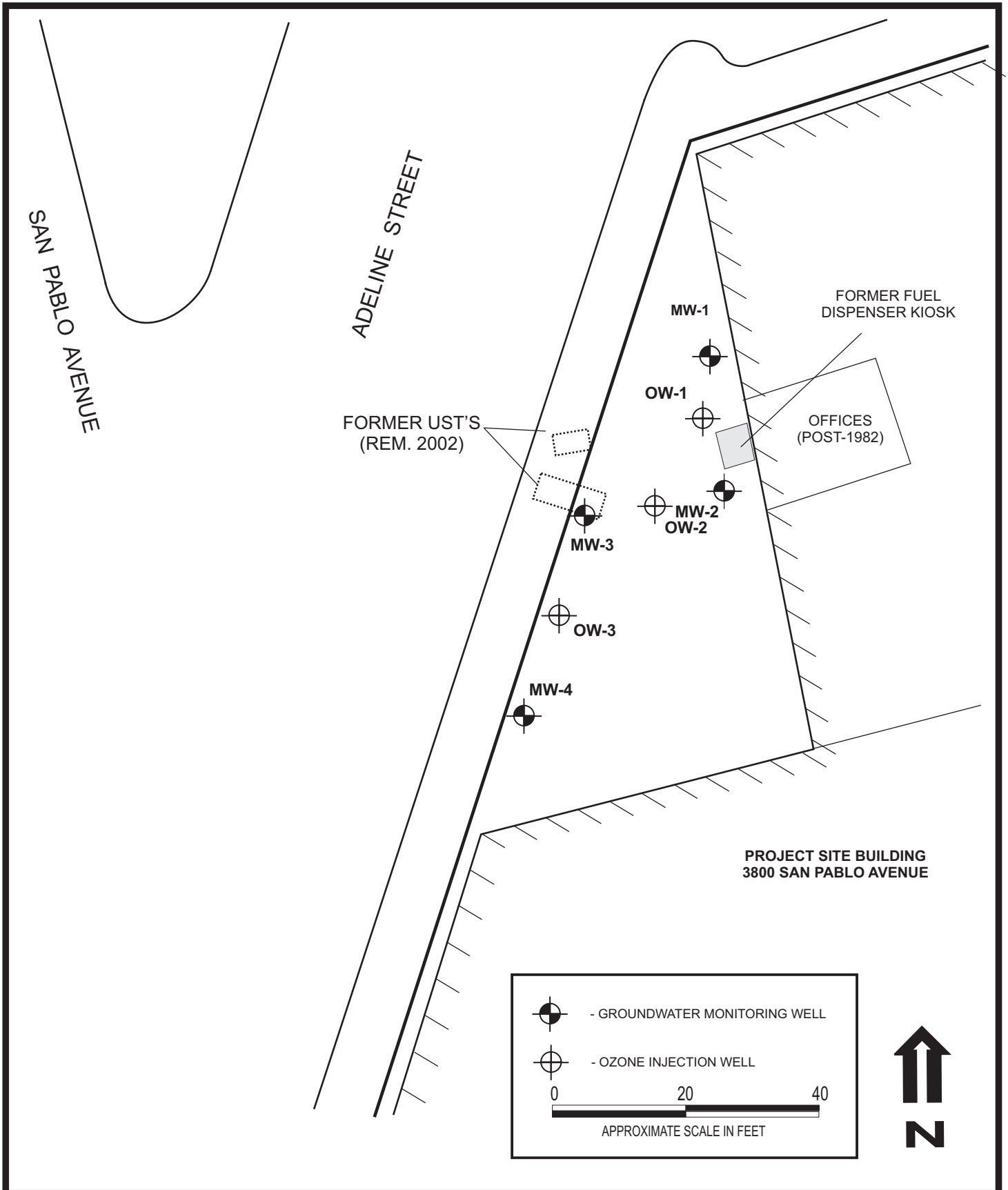
DESIGNED BY:	CHECKED BY: JG
DRAWN BY: MR	SCALE:
PROJECT NO:	

SITE VICINITY MAP

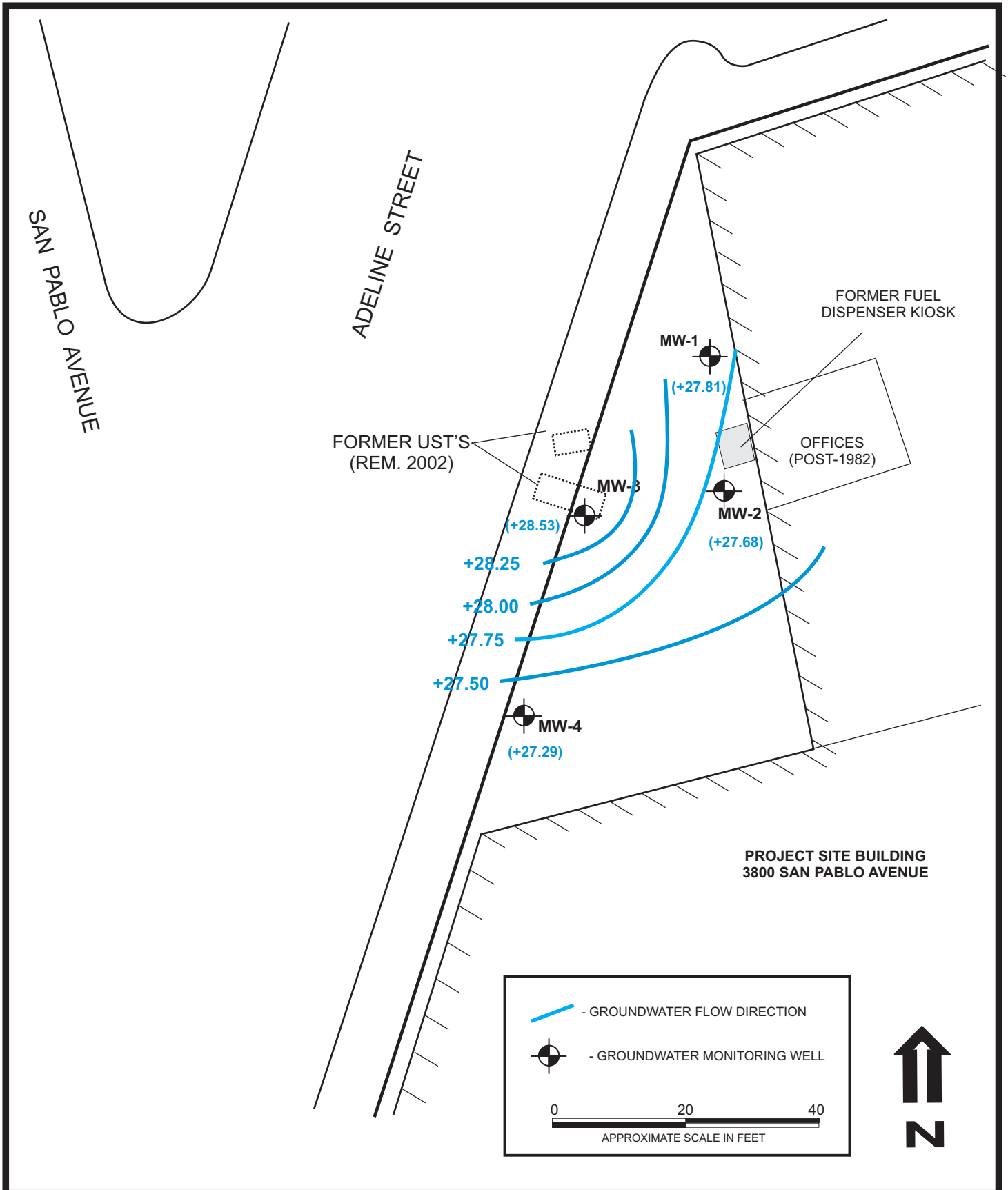
3800 SAN PABLO AVENUE
 EMERYVILLE, CALIFORNIA

DATE: 10/09/2014 FIGURE: 1

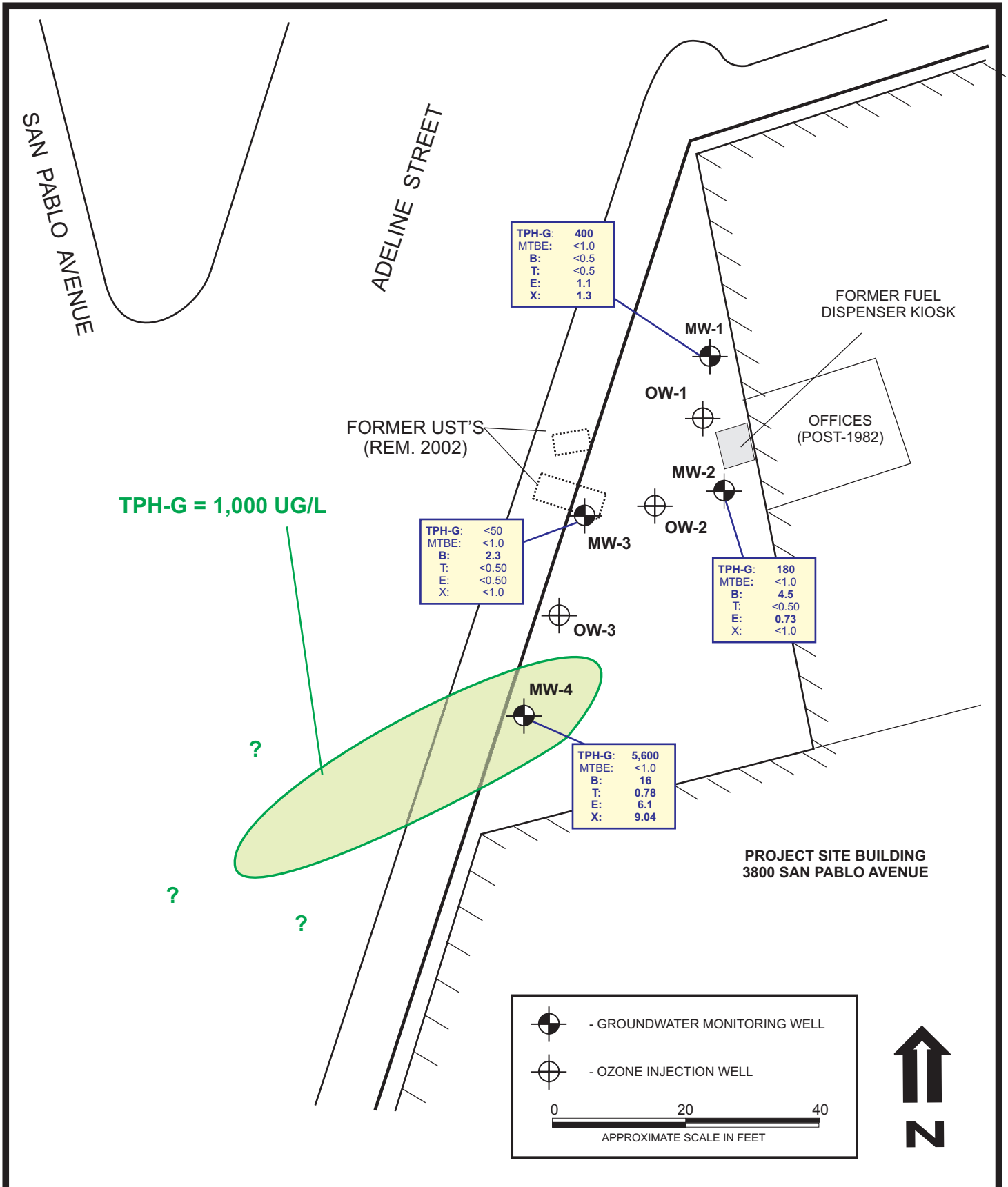




DESIGNED BY:	CHECKED BY: JG	SITE PLAN	DATE: 10/09/2014	FIGURE: 2
DRAWN BY: MR	SCALE:		GRIBI	
PROJECT NO:		3800 SAN PABLO AVENUE EMERYVILLE, CALIFORNIA		



DESIGNED BY:	CHECKED BY: JG	GROUNDWATER ELEVATION GRADIENT - 09/29/2014	DATE: 10/09/2014	FIGURE: 3
DRAWN BY: MR	SCALE:			
PROJECT NO:		3800 SAN PABLO AVENUE EMERYVILLE, CALIFORNIA		



DESIGNED BY:	CHECKED BY: JG	GROUNDWATER HYDROCARBON CONCENTRATIONS - 09/29/2014 3800 SAN PABLO AVENUE EMERYVILLE, CALIFORNIA	DATE: 10/09/2014	FIGURE: 4
DRAWN BY: MR	SCALE:			
PROJECT NO:				

ATTACHMENT A
GROUNDWATER MONITORING
FIELD DATA RECORDS

Groundwater Monitoring Field Sheet

Client Name SAN PABLO AVENUE
VENTURE Project Name MAZ GLASS
 Sampling Personnel MAR Date 9/29/2014
 Weather Conditions PC, Cool

Well ID MW-1
 Casing Diameter (inches) 2.0 Total Depth (feet) 22.7
 Depth to Water 11.75 Depth to Free Product —
 Water Column (ft) 11.45 Product Thickness φ
 One Well Volume (gal) 1.95 3x Well Volume (gal) 5.8

Notes:
 One Well Volume is determine by multiplying "Water Column" by:
 • 0.059 for 3/4-inch well, 0.17 for 2-inch well, 0.38 for 3-inch well, 0.66 for 4-inch well, 1.50 for 6-inch well

FIELD METHODS

Activity	Bailer	Pump	Comments
Purge Method		X	120 purge pump
Sample Method		X	120 purge pump

FIELD PARAMETERS

Time	Volume Purged	Temp. (F or C)	E.C. (μS/cm)	D.O. (mg/L)	pH	ORP (mV)	Comments
1228							
1230	2	21.0	856	/	7.74	/	
1233	4	20.5	903	/	7.17	/	
1235	6	20.2	895	/	7.23	/	

SAMPLE OBSERVATIONS

Characteristic	None	Slight	Moderate	Strong	Comments
Color		X			brown
Odor		X			
Turbidity		X			
Sheen	X				
Other:					

Sample Time 1235 Sampler's Signature MAR

Groundwater Monitoring Field Sheet

Client Name SAN PABLO AVENUE
VENTURE Project Name MAZ GLASS
 Sampling Personnel MAR Date 9/29/2014
 Weather Conditions PC, Cool

Well ID MW-2
 Casing Diameter (inches) 2.0 Total Depth (feet) 22.8
 Depth to Water 11.28 Depth to Free Product —
 Water Column (ft) 11.52 Product Thickness φ
 One Well Volume (gal) 1.96 3x Well Volume (gal) 5.9

Notes:
 One Well Volume is determine by multiplying "Water Column" by:
 • 0.059 for 3/4-inch well, 0.17 for 2-inch well, 0.38 for 3-inch well, 0.66 for 4-inch well, 1.50 for 6-inch well

FIELD METHODS

Activity	Bailer	Pump	Comments
Purge Method		X	120 purge pump
Sample Method		X	120 purge pump

FIELD PARAMETERS

Time	Volume Purged	Temp. (F or C)	E.C. (μS/cm)	D.O. (mg/L)	pH	ORP (mV)	Comments
1203							
1206	2	20.3	964	/	7.04	/	
1209	4	20.3	982	/	7.01	/	
1212	6	20.1	963	/	7.11	/	

SAMPLE OBSERVATIONS

Characteristic	None	Slight	Moderate	Strong	Comments
Color	X				
Odor		X →			
Turbidity	X				
Sheen	X				
Other:					

Sample Time 1215 Sampler's Signature MAR

Groundwater Monitoring Field Sheet

Client Name SAN PABLO AVENUE VENTURE Project Name MAZ GLASS
 Sampling Personnel MAR Date 9/29/2014
 Weather Conditions Cloudy, cool

Well ID MW-3
 Casing Diameter (inches) 2.0 Total Depth (feet) 22.8
 Depth to Water 10.31 Depth to Free Product
 Water Column (ft) 12.49 Product Thickness
 One Well Volume (gal) 2.12 3x Well Volume (gal) 6.4

Notes:
 One Well Volume is determine by multiplying "Water Column" by:
 • 0.059 for 3/4-inch well, 0.17 for 2-inch well, 0.38 for 3-inch well, 0.66 for 4-inch well, 1.50 for 6-inch well

FIELD METHODS

Activity	Bailer	Pump	Comments
Purge Method		X	12V purge pump
Sample Method		X	12V purge pump

FIELD PARAMETERS

Time	Volume Purged	Temp. (F or C)	E.C. (µS/cm)	D.O. (mg/L)	pH	ORP (mV)	Comments
1139							
1141	2	20.9	815	/	7.78	/	
1144	4	21.0	857	/	7.32	/	
1147	6	20.2	846	/	7.43	/	
1148	7	20.1	849	/	7.47	/	

SAMPLE OBSERVATIONS

Characteristic	None	Slight	Moderate	Strong	Comments
Color		X			brown
Odor	X				
Turbidity		X			
Sheen	X				
Other:					

Sample Time 1150 Sampler's Signature MAR

Groundwater Monitoring Field Sheet

Client Name SAN PABLO AVENUE VENTURE Project Name MAZ GLASS
 Sampling Personnel MAR Date 9/29/2014
 Weather Conditions PC, cool

Well ID MW-4
 Casing Diameter (inches) 2.0 Total Depth (feet) 22.8
 Depth to Water 11.19 Depth to Free Product
 Water Column (ft) 11.61 Product Thickness
 One Well Volume (gal) 1.97 3x Well Volume (gal) 5.9

Notes:
 One Well Volume is determine by multiplying "Water Column" by:
 • 0.059 for 3/4-inch well, 0.17 for 2-inch well, 0.38 for 3-inch well, 0.66 for 4-inch well, 1.50 for 6-inch well

FIELD METHODS

Activity	Bailer	Pump	Comments
Purge Method		X	12V purge pump
Sample Method		X	12V purge pump

FIELD PARAMETERS

Time	Volume Purged	Temp. (F or C)	E.C. (µS/cm)	D.O. (mg/L)	pH	ORP (mV)	Comments
1252							
1255	2	20.2	970	/	6.88	/	
1258	4	20.1	977	/	6.83	/	
1300	6	19.8	988	/	6.82	/	

SAMPLE OBSERVATIONS

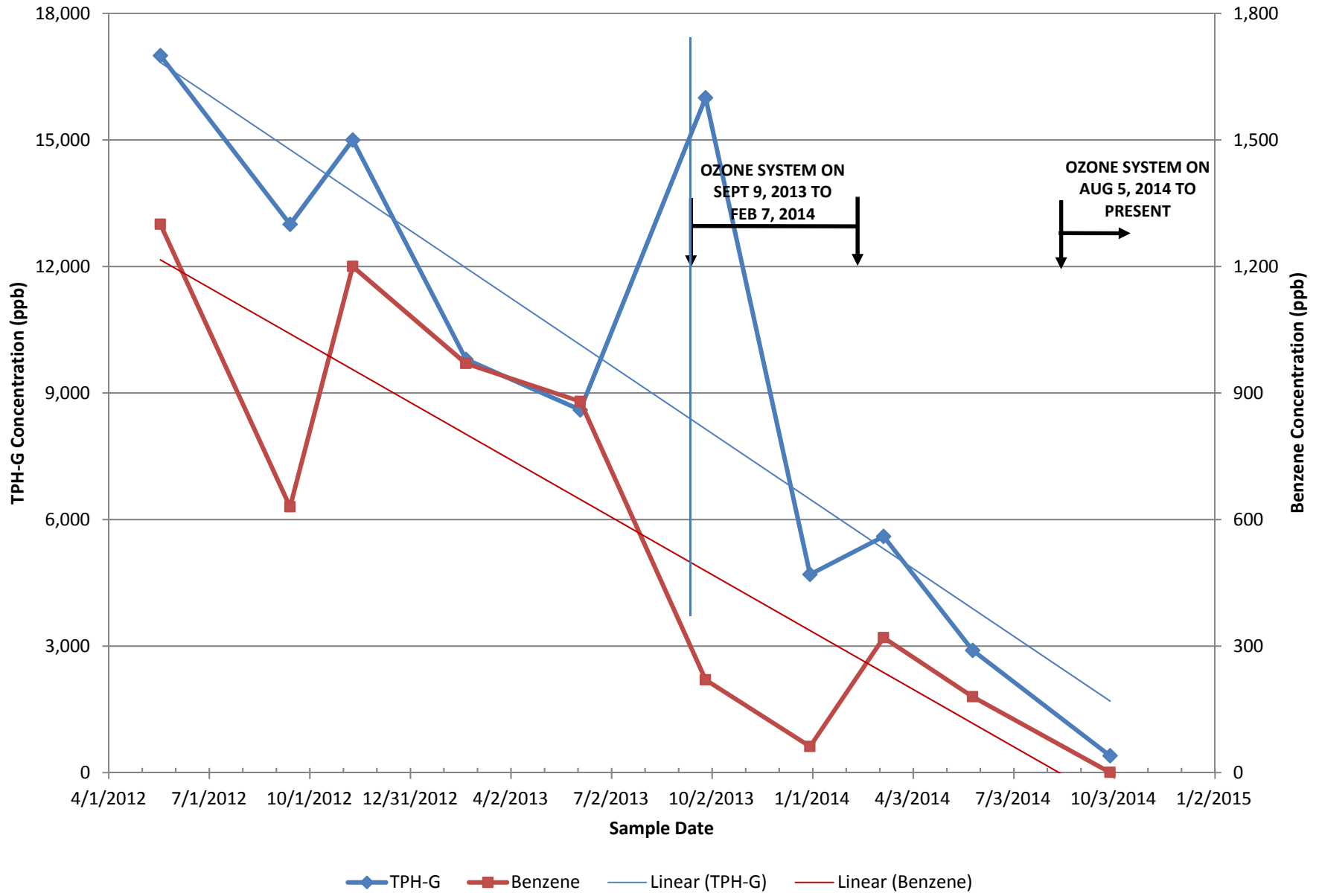
Characteristic	None	Slight	Moderate	Strong	Comments
Color	X				
Odor		X			
Turbidity	X				
Sheen	X				
Other:					

Sample Time 1300 Sampler's Signature MAR

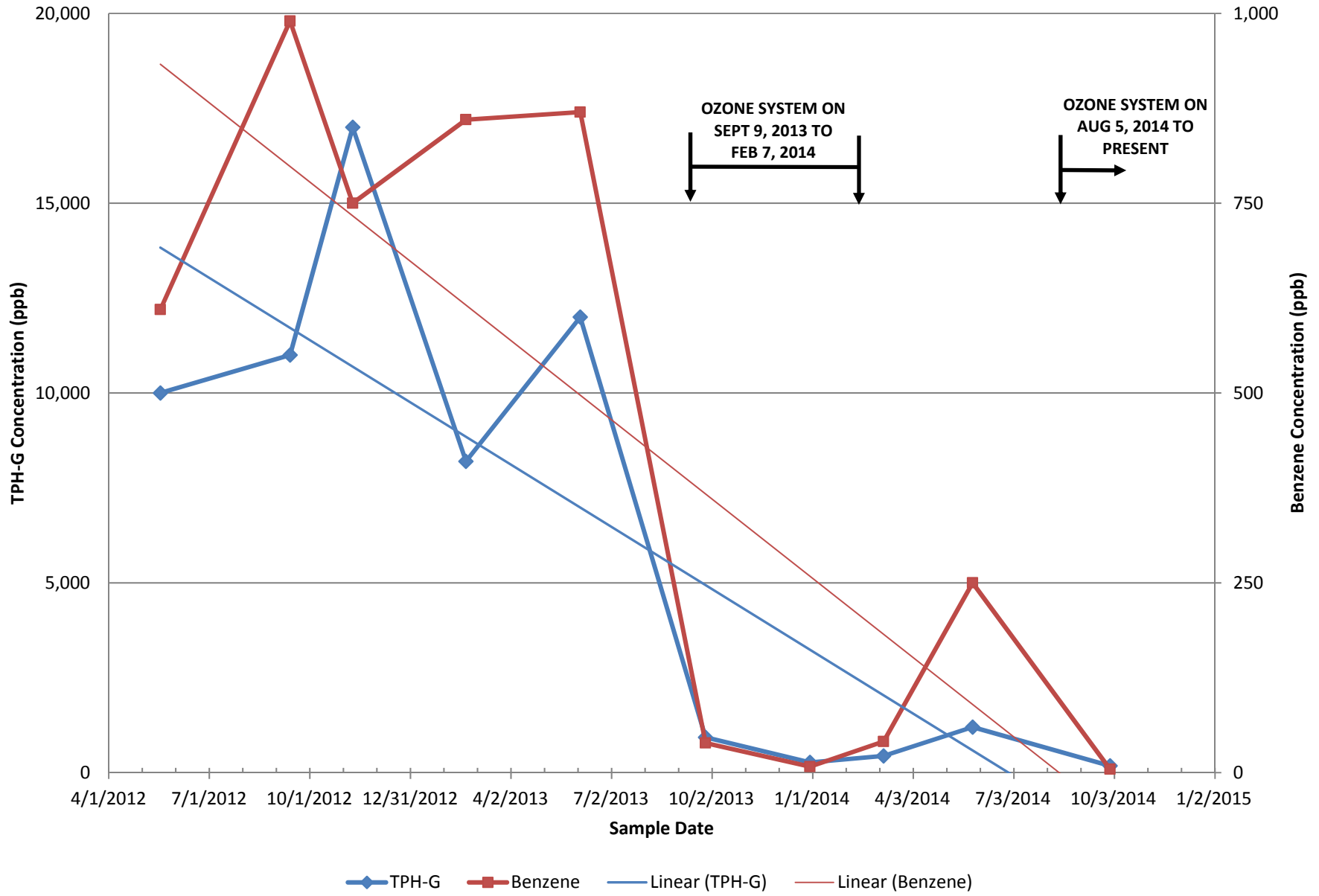
ATTACHMENT B

GROUNDWATER HYDROCARBON TRENDS

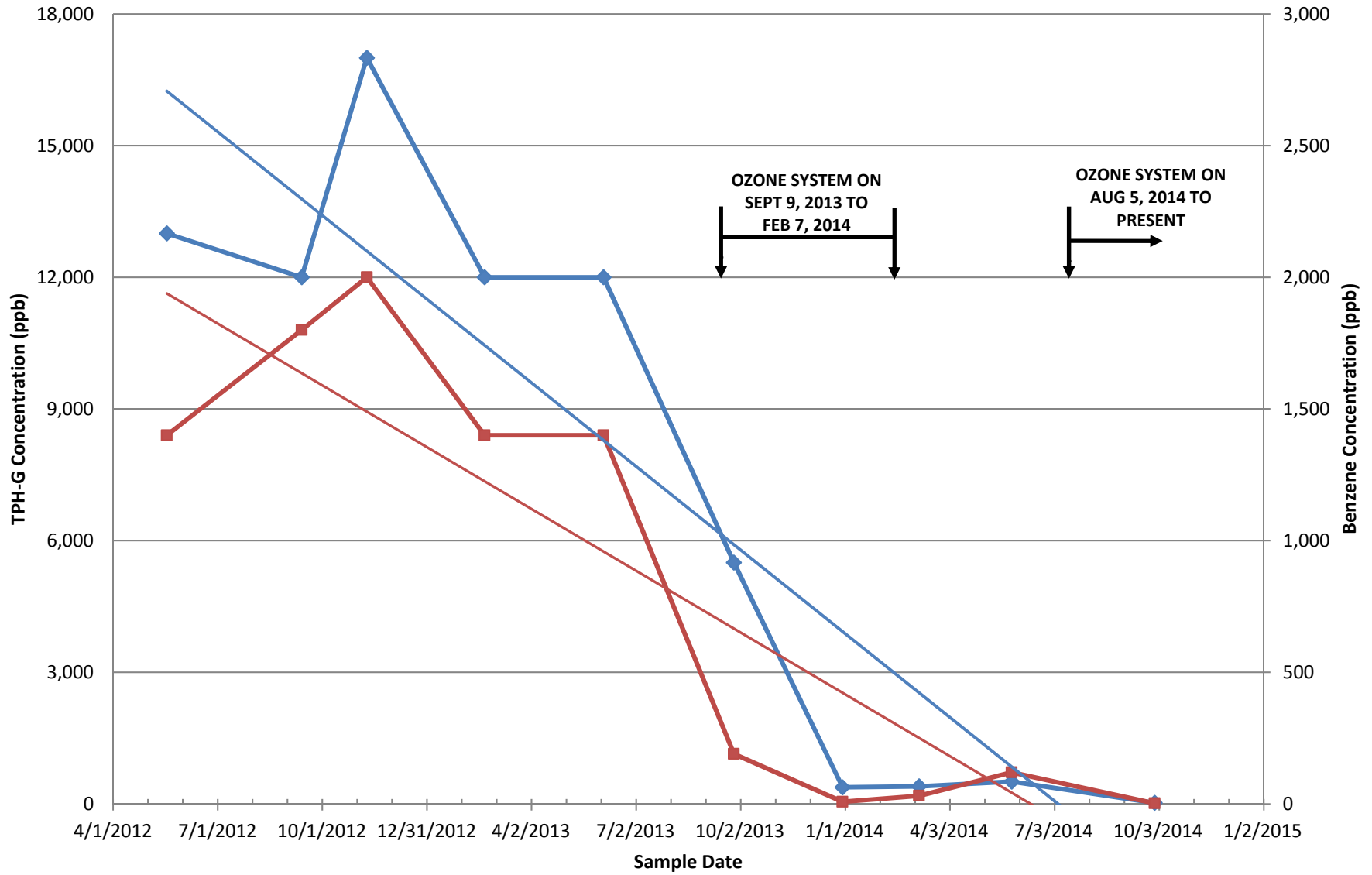
MW-1. TPH-G and Benzene Concentrations Versus Time



MW-2. TPH-G and Benzene Concentration Versus Time

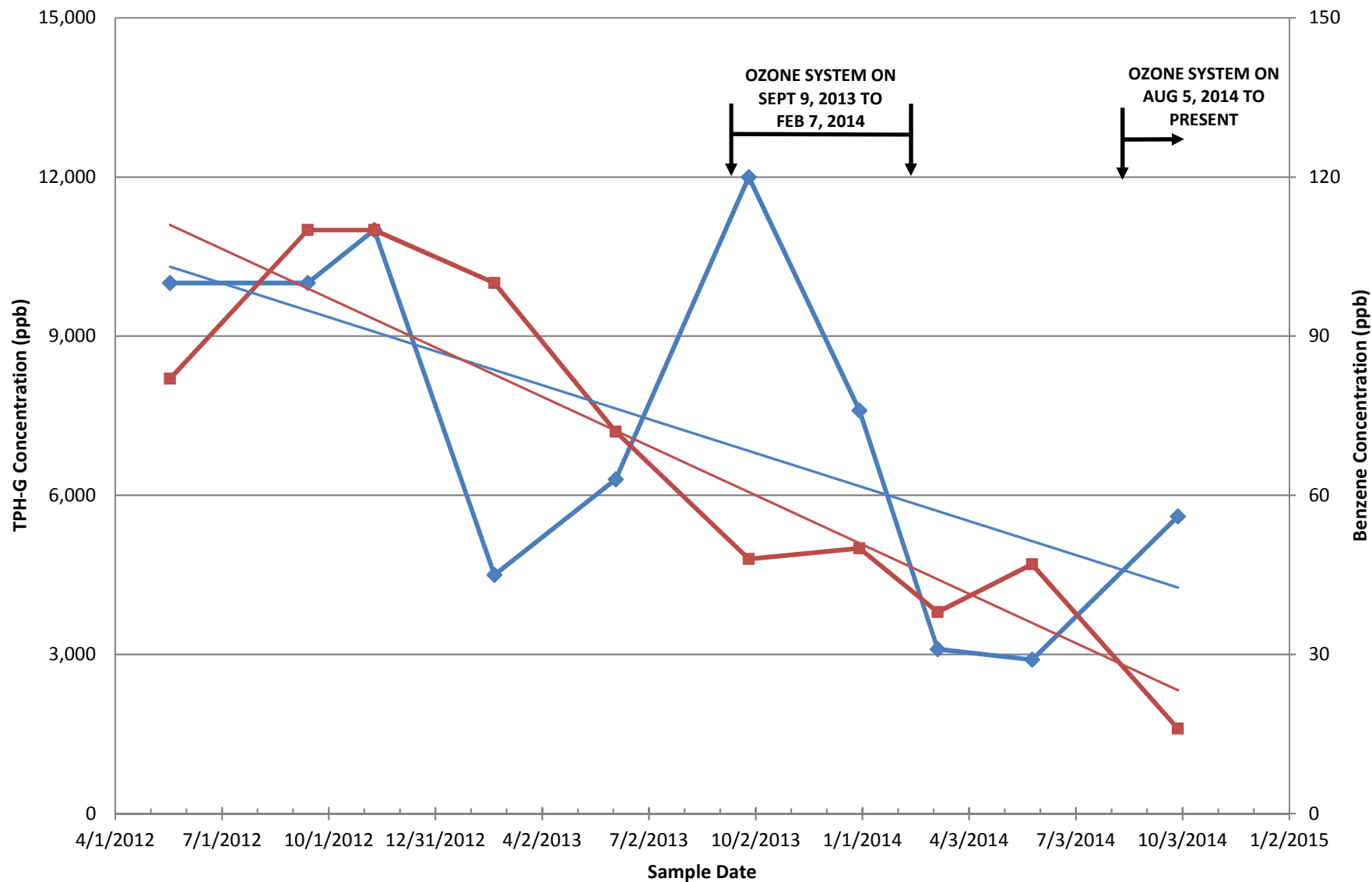


MW-3. TPH-G and Benzene Concentrations Versus Time



◆ TPH-G
 ■ Benzene
 — Linear (TPH-G)
 — Linear (Benzene)

MW-4. TPH-G and Benzene Concentrations Versus Time



◆ TPH-G
 ■ Benzene
 — Linear (TPH-G)
 — Linear (Benzene)

ATTACHMENT C

**LABORATORY DATA REPORTS AND
CHAIN-OF-CUSTODY RECORDS**



25712 Commercentre Drive
 Lake Forest, California 92630
 949.297.5020 Phone
 949.297.5027 Fax



25712 Commercentre Drive
 Lake Forest, California 92630
 949.297.5020 Phone
 949.297.5027 Fax

Gribi Associates 1090 Adam Street, Suite K Benicia CA, 94510	Project: Maz Glass Project Number: [none] Project Manager: Jim Gribi	Reported: 10/08/14 17:03
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ANALYTICAL REPORT FOR SAMPLES

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
MW-1	T142007-01	Water	09/29/14 12:35	10/01/14 08:50
MW-2	T142007-02	Water	09/29/14 12:15	10/01/14 08:50
MW-3	T142007-03	Water	09/29/14 11:50	10/01/14 08:50
MW-4	T142007-04	Water	09/29/14 13:00	10/01/14 08:50

DETECTIONS SUMMARY

08 October 2014

Jim Gribi
 Gribi Associates
 1090 Adam Street, Suite K
 Benicia, CA 94510
 RE: Maz Glass

Enclosed are the results of analyses for samples received by the laboratory on 10/01/14 08:50. If you have any questions concerning this report, please feel free to contact me.

Sample ID: MW-1		Laboratory ID: T142007-01				
Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Kerosene	0.96	0.50		mg/l	EPA 8015C	
1,3,5-Trimethylbenzene	7.0	1.0		ug/l	EPA 8260B	
1,2,4-Trimethylbenzene	4.3	1.0		ug/l	EPA 8260B	
Ethylbenzene	1.1	0.50		ug/l	EPA 8260B	
m,p-Xylene	1.3	1.0		ug/l	EPA 8260B	
Tert-butyl alcohol	38	10		ug/l	EPA 8260B	
C6-C12 (GRO)	400	50		ug/l	EPA 8260B	

Sincerely,

Katherine RunningCrane
 Project Manager

Sample ID: MW-2		Laboratory ID: T142007-02				
Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Benzene	4.5	0.50		ug/l	EPA 8260B	
Ethylbenzene	0.73	0.50		ug/l	EPA 8260B	
Tert-butyl alcohol	87	10		ug/l	EPA 8260B	
C6-C12 (GRO)	180	50		ug/l	EPA 8260B	

Sample ID: MW-3		Laboratory ID: T142007-03				
Analyte	Result	Reporting		Units	Method	Notes
		Limit				
Benzene	2.3	0.50		ug/l	EPA 8260B	

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Katherine RunningCrane, Project Manager



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Gribi Associates 1090 Adam Street, Suite K Benicia CA, 94510	Project: Maz Glass Project Number: [none] Project Manager: Jim Gribi	Reported: 10/08/14 17:03
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Sample ID: MW-4 Laboratory ID: T142007-04

Analyte	Reporting		Units	Method	Notes
	Result	Limit			
Kerosene	4.9	0.50	mg/l	EPA 8015C	
C13-C28 (DRO)	2.2	0.50	mg/l	EPA 8015C	
sec-Butylbenzene	1.3	1.0	ug/l	EPA 8260B	
Isopropylbenzene	2.8	1.0	ug/l	EPA 8260B	
p-Isopropyltoluene	2.9	1.0	ug/l	EPA 8260B	
n-Propylbenzene	5.7	1.0	ug/l	EPA 8260B	
1,3,5-Trimethylbenzene	22	1.0	ug/l	EPA 8260B	
1,2,4-Trimethylbenzene	20	1.0	ug/l	EPA 8260B	
Benzene	16	0.50	ug/l	EPA 8260B	
Toluene	0.78	0.50	ug/l	EPA 8260B	
Ethylbenzene	6.1	0.50	ug/l	EPA 8260B	
m,p-Xylene	8.5	1.0	ug/l	EPA 8260B	
o-Xylene	0.54	0.50	ug/l	EPA 8260B	
C6-C12 (GRO)	5600	50	ug/l	EPA 8260B	

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MW-1
T142007-01 (Water)

Analyte	Result	Reporting		Dilution	Batch	Prepared	Analyzed	Method	Notes
		Limit	Units						

SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015C

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Kerosene	0.96	0.50	mg/l	1	4100138	10/01/14	10/01/14	EPA 8015C	
C13-C28 (DRO)	ND	0.50	"	"	"	"	"	"	
Surrogate: p-Terphenyl	94.0 %	65-135	"	"	"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Bromobenzene	ND	1.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	5.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

MW-1
T142007-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
trans-1,2-Dichloroethene	ND	1.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	7.0	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	4.3	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	ND	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	1.1	0.50	"	"	"	"	"	"	
m,p-Xylene	1.3	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	

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Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

MW-1
T142007-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Tert-amyl methyl ether	ND	2.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
Tert-butyl alcohol	38	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
C6-C12 (GRO)	400	50	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene	101 %	83.5-119	"	"	"	"	"	"	
Surrogate: Dibromofluoromethane	111 %	81-136	"	"	"	"	"	"	
Surrogate: Toluene-d8	98.6 %	88.8-117	"	"	"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Carbazole	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
Aniline	ND	10	"	"	"	"	"	"	
Phenol	ND	10	"	"	"	"	"	"	
Acenaphthylene	ND	10	"	"	"	"	"	"	
2-Chlorophenol	ND	10	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	10	"	"	"	"	"	"	
Anthracene	ND	10	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	5.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	5.0	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	10	"	"	"	"	"	"	
2-Methylnaphthalene	ND	20	"	"	"	"	"	"	
1-Methylnaphthalene	ND	10	"	"	"	"	"	"	
Acenaphthene	ND	10	"	"	"	"	"	"	
Benzo (a) anthracene	ND	10	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	10	"	"	"	"	"	"	
4-Nitrophenol	ND	10	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	10	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	10	"	"	"	"	"	"	
Pentachlorophenol	ND	10	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	20	"	"	"	"	"	"	
Pyrene	ND	10	"	"	"	"	"	"	

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Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

MW-1
T142007-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Semivolatle Organic Compounds by EPA Method 8270C

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Benzo (a) pyrene	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
Benzyl alcohol	ND	50	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	10	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	5.0	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	20	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	10	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	5.0	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	10	"	"	"	"	"	"	
4-Chloroaniline	ND	20	"	"	"	"	"	"	
2-Chloronaphthalene	ND	10	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	20	"	"	"	"	"	"	
Chrysene	ND	10	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	10	"	"	"	"	"	"	
Dibenzofuran	ND	20	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	5.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	5.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	5.0	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	10	"	"	"	"	"	"	
Diethyl phthalate	ND	10	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	5.0	"	"	"	"	"	"	
Dimethyl phthalate	ND	10	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	5.0	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	10	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	20	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	10	"	"	"	"	"	"	
Fluoranthene	ND	5.0	"	"	"	"	"	"	
Fluorene	ND	10	"	"	"	"	"	"	
Hexachlorobenzene	ND	20	"	"	"	"	"	"	
Hexachlorobutadiene	ND	10	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	20	"	"	"	"	"	"	
Hexachloroethane	ND	5.0	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	10	"	"	"	"	"	"	

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10/08/14 17:03

MW-1
T142007-01 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Semivolatle Organic Compounds by EPA Method 8270C

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Isophorone	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
2-Methylphenol	ND	10	"	"	"	"	"	"	
4-Methylphenol	ND	20	"	"	"	"	"	"	
Naphthalene	ND	5.0	"	"	"	"	"	"	
2-Nitroaniline	ND	10	"	"	"	"	"	"	
3-Nitroaniline	ND	10	"	"	"	"	"	"	
4-Nitroaniline	ND	20	"	"	"	"	"	"	
Nitrobenzene	ND	20	"	"	"	"	"	"	
2-Nitrophenol	ND	10	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	10	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	25	"	"	"	"	"	"	
Phenanthrene	ND	10	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	20	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	10	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	10	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	10	"	"	"	"	"	"	
1,4-Dinitrobenzene	ND	10	"	"	"	"	"	"	
Pyridine	ND	10	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol	54.0 %	9.97-110	"	"	"	"	"	"	
Surrogate: Phenol-d6	48.3 %	8.4-110	"	"	"	"	"	"	
Surrogate: Nitrobenzene-d5	56.7 %	14.7-110	"	"	"	"	"	"	
Surrogate: 2-Fluorobiphenyl	52.7 %	33.3-110	"	"	"	"	"	"	
Surrogate: 2,4,6-Tribromophenol	57.1 %	12.9-110	"	"	"	"	"	"	
Surrogate: Terphenyl-d14	57.8 %	15.8-136	"	"	"	"	"	"	

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Benicia CA, 94510 Project Manager: Jim Gribi 10/08/14 17:03

MW-2
T142007-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015C

Kerosene	ND	0.50	mg/l	1	4100138	10/01/14	10/01/14	EPA 8015C	
C13-C28 (DRO)	ND	0.50	"	"	"	"	"	"	
Surrogate: p-Terphenyl		92.8 %	65-135	"	"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	5.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

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Benicia CA, 94510 Project Manager: Jim Gribi 10/08/14 17:03

MW-2
T142007-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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Volatile Organic Compounds by EPA Method 8260B

trans-1,2-Dichloroethene	ND	1.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	4.5	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	0.73	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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Gribi Associates Project: Maz Glass
1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

MW-2
T142007-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Tert-amyl methyl ether	ND	2.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
Tert-butyl alcohol	87	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
C6-C12 (GRO)	180	50	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene	101 %	83.5-119	"	"	"	"	"	"	
Surrogate: Dibromofluoromethane	114 %	81-136	"	"	"	"	"	"	
Surrogate: Toluene-d8	97.0 %	88.8-117	"	"	"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
Aniline	ND	10	"	"	"	"	"	"	
Phenol	ND	10	"	"	"	"	"	"	
Acenaphthylene	ND	10	"	"	"	"	"	"	
2-Chlorophenol	ND	10	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	10	"	"	"	"	"	"	
Anthracene	ND	10	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	5.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	5.0	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	10	"	"	"	"	"	"	
1-Methylnaphthalene	ND	10	"	"	"	"	"	"	
2-Methylnaphthalene	ND	20	"	"	"	"	"	"	
Acenaphthene	ND	10	"	"	"	"	"	"	
Benzo (a) anthracene	ND	10	"	"	"	"	"	"	
4-Nitrophenol	ND	10	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	10	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	10	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	10	"	"	"	"	"	"	
Pentachlorophenol	ND	10	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	20	"	"	"	"	"	"	
Benzo (a) pyrene	ND	10	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

MW-2
T142007-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Pyrene	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
Benzyl alcohol	ND	50	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	10	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	5.0	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	20	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	10	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	5.0	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	10	"	"	"	"	"	"	
4-Chloroaniline	ND	20	"	"	"	"	"	"	
2-Chloronaphthalene	ND	10	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	20	"	"	"	"	"	"	
Chrysene	ND	10	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	10	"	"	"	"	"	"	
Dibenzofuran	ND	20	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	5.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	5.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	5.0	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	10	"	"	"	"	"	"	
Diethyl phthalate	ND	10	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	5.0	"	"	"	"	"	"	
Dimethyl phthalate	ND	10	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	5.0	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	10	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	20	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	10	"	"	"	"	"	"	
Fluoranthene	ND	5.0	"	"	"	"	"	"	
Fluorene	ND	10	"	"	"	"	"	"	
Hexachlorobenzene	ND	20	"	"	"	"	"	"	
Hexachlorobutadiene	ND	10	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	20	"	"	"	"	"	"	
Hexachloroethane	ND	5.0	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	10	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

MW-2
T142007-02 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Isophorone	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
2-Methylphenol	ND	10	"	"	"	"	"	"	
4-Methylphenol	ND	20	"	"	"	"	"	"	
Naphthalene	ND	5.0	"	"	"	"	"	"	
2-Nitroaniline	ND	10	"	"	"	"	"	"	
3-Nitroaniline	ND	10	"	"	"	"	"	"	
4-Nitroaniline	ND	20	"	"	"	"	"	"	
Nitrobenzene	ND	20	"	"	"	"	"	"	
2-Nitrophenol	ND	10	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	10	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	25	"	"	"	"	"	"	
Phenanthrene	ND	10	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	20	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	10	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	10	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	10	"	"	"	"	"	"	
1,4-Dinitrobenzene	ND	10	"	"	"	"	"	"	
Pyridine	ND	10	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol	65.6 %	9.97-110	"	"	"	"	"	"	
Surrogate: Phenol-d6	54.2 %	8.4-110	"	"	"	"	"	"	
Surrogate: Nitrobenzene-d5	56.2 %	14.7-110	"	"	"	"	"	"	
Surrogate: 2-Fluorobiphenyl	43.3 %	33.3-110	"	"	"	"	"	"	
Surrogate: 2,4,6-Tribromophenol	67.6 %	12.9-110	"	"	"	"	"	"	
Surrogate: Terphenyl-d14	42.2 %	15.8-136	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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Gribi Associates Project: Maz Glass
1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

MW-3
T142007-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015C

Kerosene	ND	0.50	mg/l	1	4100138	10/01/14	10/01/14	EPA 8015C	
C13-C28 (DRO)	ND	0.50	"	"	"	"	"	"	
Surrogate: p-Terphenyl	97.5 %	65-135	"	"	"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	ND	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	5.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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Gribi Associates Project: Maz Glass
1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

MW-3
T142007-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
trans-1,2-Dichloroethene	ND	1.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
1,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	ND	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	ND	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	ND	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	ND	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	2.3	0.50	"	"	"	"	"	"	
Toluene	ND	0.50	"	"	"	"	"	"	
Ethylbenzene	ND	0.50	"	"	"	"	"	"	
m,p-Xylene	ND	1.0	"	"	"	"	"	"	
o-Xylene	ND	0.50	"	"	"	"	"	"	

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Katherine RunningCrane

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1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

MW-3
T142007-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Tert-amyl methyl ether	ND	2.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	
Di-isopropyl ether	ND	2.0	"	"	"	"	"	"	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
C6-C12 (GRO)	ND	50	"	"	"	"	"	"	
Surrogate: 4-Bromofluorobenzene	95.4 %	83.5-119	"	"	"	"	"	"	
Surrogate: Dibromofluoromethane	112 %	81-136	"	"	"	"	"	"	
Surrogate: Toluene-d8	96.5 %	88.8-117	"	"	"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Carbazole	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
Aniline	ND	10	"	"	"	"	"	"	
Phenol	ND	10	"	"	"	"	"	"	
Acenaphthylene	ND	10	"	"	"	"	"	"	
2-Chlorophenol	ND	10	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	10	"	"	"	"	"	"	
Anthracene	ND	10	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	5.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	5.0	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	10	"	"	"	"	"	"	
1-Methylnaphthalene	ND	10	"	"	"	"	"	"	
2-Methylnaphthalene	ND	20	"	"	"	"	"	"	
Acenaphthene	ND	10	"	"	"	"	"	"	
Benzo (a) anthracene	ND	10	"	"	"	"	"	"	
4-Nitrophenol	ND	10	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	10	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	10	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	10	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	20	"	"	"	"	"	"	
Pentachlorophenol	ND	10	"	"	"	"	"	"	
Benzo (a) pyrene	ND	10	"	"	"	"	"	"	

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Gribi Associates Project: Maz Glass
 1090 Adam Street, Suite K Project Number: [none]
 Benicia CA, 94510 Project Manager: Jim Gribi Reported:
 10/08/14 17:03

MW-3
T142007-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Pyrene	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
Benzyl alcohol	ND	50	"	"	"	"	"	"	
Bis(2-chloroethoxy)methane	ND	10	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	5.0	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	20	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	10	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	5.0	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	10	"	"	"	"	"	"	
4-Chloroaniline	ND	20	"	"	"	"	"	"	
2-Chloronaphthalene	ND	10	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	20	"	"	"	"	"	"	
Chrysene	ND	10	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	10	"	"	"	"	"	"	
Dibenzofuran	ND	20	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	5.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	5.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	5.0	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	10	"	"	"	"	"	"	
Diethyl phthalate	ND	10	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	5.0	"	"	"	"	"	"	
Dimethyl phthalate	ND	10	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	5.0	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	10	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	20	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	10	"	"	"	"	"	"	
Fluoranthene	ND	5.0	"	"	"	"	"	"	
Fluorene	ND	10	"	"	"	"	"	"	
Hexachlorobenzene	ND	20	"	"	"	"	"	"	
Hexachlorobutadiene	ND	10	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	20	"	"	"	"	"	"	
Hexachloroethane	ND	5.0	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	10	"	"	"	"	"	"	

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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Gribi Associates Project: Maz Glass
 1090 Adam Street, Suite K Project Number: [none]
 Benicia CA, 94510 Project Manager: Jim Gribi Reported:
 10/08/14 17:03

MW-3
T142007-03 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
Isophorone	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
2-Methylphenol	ND	10	"	"	"	"	"	"	
4-Methylphenol	ND	20	"	"	"	"	"	"	
Naphthalene	ND	5.0	"	"	"	"	"	"	
2-Nitroaniline	ND	10	"	"	"	"	"	"	
3-Nitroaniline	ND	10	"	"	"	"	"	"	
4-Nitroaniline	ND	20	"	"	"	"	"	"	
Nitrobenzene	ND	20	"	"	"	"	"	"	
2-Nitrophenol	ND	10	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	10	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	25	"	"	"	"	"	"	
Phenanthrene	ND	10	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	20	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	10	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	10	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	10	"	"	"	"	"	"	
1,4-Dinitrobenzene	ND	10	"	"	"	"	"	"	
Pyridine	ND	10	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol	48.3 %	9.97-110	"	"	"	"	"	"	
Surrogate: Phenol-d6	41.8 %	8.4-110	"	"	"	"	"	"	
Surrogate: Nitrobenzene-d5	59.2 %	14.7-110	"	"	"	"	"	"	
Surrogate: 2-Fluorobiphenyl	45.8 %	33.3-110	"	"	"	"	"	"	
Surrogate: 2,4,6-Tribromophenol	60.3 %	12.9-110	"	"	"	"	"	"	
Surrogate: Terphenyl-d14	58.6 %	15.8-136	"	"	"	"	"	"	

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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 Benicia CA, 94510 Project Manager: Jim Gribi 10/08/14 17:03

MW-4
T142007-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Extractable Petroleum Hydrocarbons by 8015C

Kerosene	4.9	0.50	mg/l	1	4100138	10/01/14	10/01/14	EPA 8015C	
C13-C28 (DRO)	2.2	0.50	"	"	"	"	"	"	
<i>Surrogate: p-Terphenyl</i>		90.6 %	65-135	"	"	"	"	"	

Volatile Organic Compounds by EPA Method 8260B

Bromobenzene	ND	1.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
Bromochloromethane	ND	1.0	"	"	"	"	"	"	
Bromodichloromethane	ND	1.0	"	"	"	"	"	"	
Bromoform	ND	1.0	"	"	"	"	"	"	
Bromomethane	ND	1.0	"	"	"	"	"	"	
n-Butylbenzene	ND	1.0	"	"	"	"	"	"	
sec-Butylbenzene	1.3	1.0	"	"	"	"	"	"	
tert-Butylbenzene	ND	1.0	"	"	"	"	"	"	
Carbon tetrachloride	ND	0.50	"	"	"	"	"	"	
Chlorobenzene	ND	1.0	"	"	"	"	"	"	
Chloroethane	ND	1.0	"	"	"	"	"	"	
Chloroform	ND	1.0	"	"	"	"	"	"	
Chloromethane	ND	1.0	"	"	"	"	"	"	
2-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
4-Chlorotoluene	ND	1.0	"	"	"	"	"	"	
Dibromochloromethane	ND	1.0	"	"	"	"	"	"	
1,2-Dibromo-3-chloropropane	ND	5.0	"	"	"	"	"	"	
1,2-Dibromoethane (EDB)	ND	1.0	"	"	"	"	"	"	
Dibromomethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	1.0	"	"	"	"	"	"	
Dichlorodifluoromethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethane	ND	1.0	"	"	"	"	"	"	
1,2-Dichloroethane	ND	0.50	"	"	"	"	"	"	
1,1-Dichloroethene	ND	1.0	"	"	"	"	"	"	
cis-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	
trans-1,2-Dichloroethene	ND	1.0	"	"	"	"	"	"	

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 Benicia CA, 94510 Project Manager: Jim Gribi 10/08/14 17:03

MW-4
T142007-04 (Water)

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

1,2-Dichloropropane	ND	1.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
1,3-Dichloropropane	ND	1.0	"	"	"	"	"	"	
2,2-Dichloropropane	ND	1.0	"	"	"	"	"	"	
1,1-Dichloropropene	ND	1.0	"	"	"	"	"	"	
cis-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
trans-1,3-Dichloropropene	ND	0.50	"	"	"	"	"	"	
Hexachlorobutadiene	ND	1.0	"	"	"	"	"	"	
Isopropylbenzene	2.8	1.0	"	"	"	"	"	"	
p-Isopropyltoluene	2.9	1.0	"	"	"	"	"	"	
Methylene chloride	ND	1.0	"	"	"	"	"	"	
Naphthalene	ND	1.0	"	"	"	"	"	"	
n-Propylbenzene	5.7	1.0	"	"	"	"	"	"	
Styrene	ND	1.0	"	"	"	"	"	"	
1,1,2,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1,2-Tetrachloroethane	ND	1.0	"	"	"	"	"	"	
Tetrachloroethene	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	1.0	"	"	"	"	"	"	
1,1,2-Trichloroethane	ND	1.0	"	"	"	"	"	"	
1,1,1-Trichloroethane	ND	1.0	"	"	"	"	"	"	
Trichloroethene	ND	1.0	"	"	"	"	"	"	
Trichlorofluoromethane	ND	1.0	"	"	"	"	"	"	
1,2,3-Trichloropropane	ND	1.0	"	"	"	"	"	"	
1,3,5-Trimethylbenzene	22	1.0	"	"	"	"	"	"	
1,2,4-Trimethylbenzene	20	1.0	"	"	"	"	"	"	
Vinyl chloride	ND	1.0	"	"	"	"	"	"	
Benzene	16	0.50	"	"	"	"	"	"	
Toluene	0.78	0.50	"	"	"	"	"	"	
Ethylbenzene	6.1	0.50	"	"	"	"	"	"	
m,p-Xylene	8.5	1.0	"	"	"	"	"	"	
o-Xylene	0.54	0.50	"	"	"	"	"	"	
Tert-amyl methyl ether	ND	2.0	"	"	"	"	"	"	
Tert-butyl alcohol	ND	10	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine RunningCrane, Project Manager



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Gribi Associates Project: Maz Glass
1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

**MW-4
T142007-04 (Water)**

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Volatile Organic Compounds by EPA Method 8260B

Di-isopropyl ether	ND	2.0	ug/l	1	4100140	10/01/14	10/01/14	EPA 8260B	
Ethyl tert-butyl ether	ND	2.0	"	"	"	"	"	"	
Methyl tert-butyl ether	ND	1.0	"	"	"	"	"	"	
Ethanol	ND	500	"	"	"	"	"	"	
C6-C12 (GRO)	5600	50	"	"	"	"	"	"	"
Surrogate: 4-Bromofluorobenzene	98.4 %	83.5-119	"	"	"	"	"	"	
Surrogate: Dibromofluoromethane	107 %	81-136	"	"	"	"	"	"	
Surrogate: Toluene-d8	97.5 %	88.8-117	"	"	"	"	"	"	

Semivolatile Organic Compounds by EPA Method 8270C

Carbazole	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
Aniline	ND	10	"	"	"	"	"	"	
Phenol	ND	10	"	"	"	"	"	"	
Acenaphthylene	ND	10	"	"	"	"	"	"	
2-Chlorophenol	ND	10	"	"	"	"	"	"	
1,4-Dichlorobenzene	ND	10	"	"	"	"	"	"	
Anthracene	ND	10	"	"	"	"	"	"	
N-Nitrosodi-n-propylamine	ND	5.0	"	"	"	"	"	"	
1,2,4-Trichlorobenzene	ND	5.0	"	"	"	"	"	"	
4-Chloro-3-methylphenol	ND	10	"	"	"	"	"	"	
1-Methylnaphthalene	ND	10	"	"	"	"	"	"	
2-Methylnaphthalene	ND	20	"	"	"	"	"	"	
Benzo (a) anthracene	ND	10	"	"	"	"	"	"	
Acenaphthene	ND	10	"	"	"	"	"	"	
4-Nitrophenol	ND	10	"	"	"	"	"	"	
Benzo (b) fluoranthene	ND	10	"	"	"	"	"	"	
2,4-Dinitrotoluene	ND	10	"	"	"	"	"	"	
Benzo (k) fluoranthene	ND	10	"	"	"	"	"	"	
Pentachlorophenol	ND	10	"	"	"	"	"	"	
Benzo (g,h,i) perylene	ND	20	"	"	"	"	"	"	
Benzo (a) pyrene	ND	10	"	"	"	"	"	"	
Pyrene	ND	10	"	"	"	"	"	"	

SunStar Laboratories, Inc.

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

**MW-4
T142007-04 (Water)**

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Semivolatile Organic Compounds by EPA Method 8270C

Benzyl alcohol	ND	50	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
Bis(2-chloroethoxy)methane	ND	10	"	"	"	"	"	"	
Bis(2-chloroethyl)ether	ND	5.0	"	"	"	"	"	"	
Bis(2-chloroisopropyl)ether	ND	20	"	"	"	"	"	"	
Bis(2-ethylhexyl)phthalate	ND	10	"	"	"	"	"	"	
4-Bromophenyl phenyl ether	ND	5.0	"	"	"	"	"	"	
Butyl benzyl phthalate	ND	10	"	"	"	"	"	"	
4-Chloroaniline	ND	20	"	"	"	"	"	"	
2-Chloronaphthalene	ND	10	"	"	"	"	"	"	
4-Chlorophenyl phenyl ether	ND	20	"	"	"	"	"	"	
Chrysene	ND	10	"	"	"	"	"	"	
Dibenz (a,h) anthracene	ND	10	"	"	"	"	"	"	
Dibenzofuran	ND	20	"	"	"	"	"	"	
Di-n-butyl phthalate	ND	5.0	"	"	"	"	"	"	
1,2-Dichlorobenzene	ND	5.0	"	"	"	"	"	"	
1,3-Dichlorobenzene	ND	5.0	"	"	"	"	"	"	
2,4-Dichlorophenol	ND	10	"	"	"	"	"	"	
Diethyl phthalate	ND	10	"	"	"	"	"	"	
2,4-Dimethylphenol	ND	5.0	"	"	"	"	"	"	
Dimethyl phthalate	ND	10	"	"	"	"	"	"	
4,6-Dinitro-2-methylphenol	ND	5.0	"	"	"	"	"	"	
2,4-Dinitrophenol	ND	10	"	"	"	"	"	"	
2,6-Dinitrotoluene	ND	20	"	"	"	"	"	"	
Di-n-octyl phthalate	ND	10	"	"	"	"	"	"	
Fluoranthene	ND	5.0	"	"	"	"	"	"	
Fluorene	ND	10	"	"	"	"	"	"	
Hexachlorobenzene	ND	20	"	"	"	"	"	"	
Hexachlorobutadiene	ND	10	"	"	"	"	"	"	
Hexachlorocyclopentadiene	ND	20	"	"	"	"	"	"	
Hexachloroethane	ND	5.0	"	"	"	"	"	"	
Indeno (1,2,3-cd) pyrene	ND	10	"	"	"	"	"	"	
Isophorone	ND	10	"	"	"	"	"	"	

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**MW-4
 T142007-04 (Water)**

Analyte	Result	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed	Method	Notes
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SunStar Laboratories, Inc.

Semivolatle Organic Compounds by EPA Method 8270C

2-Methylphenol	ND	10	ug/l	1	4100222	10/02/14	10/03/14	EPA 8270C	
4-Methylphenol	ND	20	"	"	"	"	"	"	
Naphthalene	ND	5.0	"	"	"	"	"	"	
2-Nitroaniline	ND	10	"	"	"	"	"	"	
3-Nitroaniline	ND	10	"	"	"	"	"	"	
4-Nitroaniline	ND	20	"	"	"	"	"	"	
Nitrobenzene	ND	20	"	"	"	"	"	"	
2-Nitrophenol	ND	10	"	"	"	"	"	"	
N-Nitrosodiphenylamine	ND	10	"	"	"	"	"	"	
N-Nitrosodimethylamine	ND	25	"	"	"	"	"	"	
Phenanthrene	ND	10	"	"	"	"	"	"	
2,4,5-Trichlorophenol	ND	20	"	"	"	"	"	"	
2,4,6-Trichlorophenol	ND	10	"	"	"	"	"	"	
2,3,4,6-Tetrachlorophenol	ND	10	"	"	"	"	"	"	
2,3,5,6-Tetrachlorophenol	ND	10	"	"	"	"	"	"	
1,4-Dinitrobenzene	ND	10	"	"	"	"	"	"	
Pyridine	ND	10	"	"	"	"	"	"	
Surrogate: 2-Fluorophenol		47.7 %		9.97-110	"	"	"	"	
Surrogate: Phenol-d6		50.8 %		8.4-110	"	"	"	"	
Surrogate: Nitrobenzene-d5		56.2 %		14.7-110	"	"	"	"	
Surrogate: 2-Fluorobiphenyl		41.0 %		33.3-110	"	"	"	"	
Surrogate: 2,4,6-Tribromophenol		62.8 %		12.9-110	"	"	"	"	
Surrogate: Terphenyl-d14		46.0 %		15.8-136	"	"	"	"	

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**Extractable Petroleum Hydrocarbons by 8015C - Quality Control
 SunStar Laboratories, Inc.**

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 4100138 - EPA 3510C GC

Blank (4100138-BLK1)										
Prepared & Analyzed: 10/01/14										
Kerosene	ND	0.50	mg/l							
C13-C28 (DRO)	ND	0.50	"							
Surrogate: p-Terphenyl	3.44		"	4.00		86.1	65-135			
LCS (4100138-BS1)										
Prepared & Analyzed: 10/01/14										
C13-C28 (DRO)	18.1	0.50	mg/l	20.0	0.273	90.3	75-125			
Surrogate: p-Terphenyl	3.68		"	4.00		91.9	65-135			
Matrix Spike (4100138-MS1)										
Source: T142007-01 Prepared & Analyzed: 10/01/14										
C13-C28 (DRO)	18.3	0.50	mg/l	20.0	0.273	90.1	75-125			
Surrogate: p-Terphenyl	3.77		"	4.00		94.2	65-135			
Matrix Spike Dup (4100138-MSD1)										
Source: T142007-01 Prepared & Analyzed: 10/01/14										
C13-C28 (DRO)	18.5	0.50	mg/l	20.0	0.273	91.1	75-125	1.03	20	
Surrogate: p-Terphenyl	3.56		"	4.00		89.0	65-135			

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Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 4100140 - EPA 5030 GCMS

Prepared & Analyzed: 10/01/14										
Blank (4100140-BLK1)										
Bromobenzene	ND	1.0	ug/l							
Bromochloromethane	ND	1.0	"							
Bromodichloromethane	ND	1.0	"							
Bromoform	ND	1.0	"							
Bromomethane	ND	1.0	"							
n-Butylbenzene	ND	1.0	"							
sec-Butylbenzene	ND	1.0	"							
tert-Butylbenzene	ND	1.0	"							
Carbon tetrachloride	ND	0.50	"							
Chlorobenzene	ND	1.0	"							
Chloroethane	ND	1.0	"							
Chloroform	ND	1.0	"							
Chloromethane	ND	1.0	"							
2-Chlorotoluene	ND	1.0	"							
4-Chlorotoluene	ND	1.0	"							
Dibromochloromethane	ND	1.0	"							
1,2-Dibromo-3-chloropropane	ND	5.0	"							
1,2-Dibromoethane (EDB)	ND	1.0	"							
Dibromomethane	ND	1.0	"							
1,2-Dichlorobenzene	ND	1.0	"							
1,3-Dichlorobenzene	ND	1.0	"							
1,4-Dichlorobenzene	ND	1.0	"							
Dichlorodifluoromethane	ND	0.50	"							
1,1-Dichloroethane	ND	1.0	"							
1,2-Dichloroethane	ND	0.50	"							
1,1-Dichloroethene	ND	1.0	"							
cis-1,2-Dichloroethene	ND	1.0	"							
trans-1,2-Dichloroethene	ND	1.0	"							
1,2-Dichloropropane	ND	1.0	"							
1,3-Dichloropropane	ND	1.0	"							
2,2-Dichloropropane	ND	1.0	"							
1,1-Dichloropropene	ND	1.0	"							
cis-1,3-Dichloropropene	ND	0.50	"							
trans-1,3-Dichloropropene	ND	0.50	"							
Hexachlorobutadiene	ND	1.0	"							
Isopropylbenzene	ND	1.0	"							

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Katherine RunningCrane

Katherine RunningCrane, Project Manager



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Gribi Associates Project: Maz Glass
1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 4100140 - EPA 5030 GCMS

Prepared & Analyzed: 10/01/14										
Blank (4100140-BLK1)										
p-Isopropyltoluene	ND	1.0	ug/l							
Methylene chloride	ND	1.0	"							
Naphthalene	ND	1.0	"							
n-Propylbenzene	ND	1.0	"							
Styrene	ND	1.0	"							
1,1,2,2-Tetrachloroethane	ND	1.0	"							
1,1,1,2-Tetrachloroethane	ND	1.0	"							
Tetrachloroethene	ND	1.0	"							
1,2,3-Trichlorobenzene	ND	1.0	"							
1,2,4-Trichlorobenzene	ND	1.0	"							
1,1,2-Trichloroethane	ND	1.0	"							
1,1,1-Trichloroethane	ND	1.0	"							
Trichloroethene	ND	1.0	"							
Trichlorofluoromethane	ND	1.0	"							
1,2,3-Trichloropropane	ND	1.0	"							
1,3,5-Trimethylbenzene	ND	1.0	"							
1,2,4-Trimethylbenzene	ND	1.0	"							
Vinyl chloride	ND	1.0	"							
Benzene	ND	0.50	"							
Toluene	ND	0.50	"							
Ethylbenzene	ND	0.50	"							
m,p-Xylene	ND	1.0	"							
o-Xylene	ND	0.50	"							
Tert-amyl methyl ether	ND	2.0	"							
Tert-butyl alcohol	ND	10	"							
Di-isopropyl ether	ND	2.0	"							
Ethyl tert-butyl ether	ND	2.0	"							
Methyl tert-butyl ether	ND	1.0	"							
C6-C12 (GRO)	ND	50	"							
Surrogate: 4-Bromofluorobenzene	7.98	"		8.00		99.8	83.5-119			
Surrogate: Dibromofluoromethane	8.95	"		8.00		112	81-136			
Surrogate: Toluene-d8	7.89	"		8.00		98.6	88.8-117			

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Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

Volatile Organic Compounds by EPA Method 8260B - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 4100140 - EPA 5030 GCMS

LCS (4100140-BS1)		Prepared: 10/01/14 Analyzed: 10/02/14							
Chlorobenzene	21.6	1.0	ug/l	20.0	108	75-125			
1,1-Dichloroethene	16.1	1.0	"	20.0	80.6	75-125			
Trichloroethene	17.4	1.0	"	20.0	87.2	75-125			
Benzene	17.7	0.50	"	20.0	88.6	75-125			
Toluene	17.3	0.50	"	20.0	86.4	75-125			
Surrogate: 4-Bromofluorobenzene	8.12		"	8.00	102	83.5-119			
Surrogate: Dibromofluoromethane	10.0		"	8.00	126	81-136			
Surrogate: Toluene-d8	7.30		"	8.00	91.2	88.8-117			

Matrix Spike (4100140-MS1)		Source: T142007-02 Prepared: 10/01/14 Analyzed: 10/02/14							
Chlorobenzene	21.1	1.0	ug/l	20.0	ND	106	75-125		
1,1-Dichloroethene	16.3	1.0	"	20.0	ND	81.5	75-125		
Trichloroethene	16.2	1.0	"	20.0	ND	81.2	75-125		
Benzene	23.4	0.50	"	20.0	4.50	94.4	75-125		
Toluene	17.4	0.50	"	20.0	ND	87.2	75-125		
Surrogate: 4-Bromofluorobenzene	7.67		"	8.00	95.9	83.5-119			
Surrogate: Dibromofluoromethane	10.6		"	8.00	132	81-136			
Surrogate: Toluene-d8	7.35		"	8.00	91.9	88.8-117			

Matrix Spike Dup (4100140-MSD1)		Source: T142007-02 Prepared: 10/01/14 Analyzed: 10/02/14							
Chlorobenzene	21.8	1.0	ug/l	20.0	ND	109	75-125	3.21	20
1,1-Dichloroethene	16.2	1.0	"	20.0	ND	81.2	75-125	0.307	20
Trichloroethene	16.1	1.0	"	20.0	ND	80.4	75-125	0.990	20
Benzene	23.9	0.50	"	20.0	4.50	96.8	75-125	2.03	20
Toluene	16.8	0.50	"	20.0	ND	83.8	75-125	3.86	20
Surrogate: 4-Bromofluorobenzene	8.56		"	8.00	107	83.5-119			
Surrogate: Dibromofluoromethane	10.0		"	8.00	125	81-136			
Surrogate: Toluene-d8	7.41		"	8.00	92.6	88.8-117			

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Katherine RunningCrane, Project Manager



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Gribi Associates Project: Maz Glass
1090 Adam Street, Suite K Project Number: [none]
Benicia CA, 94510 Project Manager: Jim Gribi Reported:
10/08/14 17:03

Semivolatile Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 4100222 - EPA 3510C GCMS/ECD

Blank (4100222-BLK1)		Prepared: 10/02/14 Analyzed: 10/03/14							
Carbazole	ND	10	ug/l						
Phenol	ND	10	"						
Aniline	ND	10	"						
Acenaphthylene	ND	10	"						
2-Chlorophenol	ND	10	"						
1,4-Dichlorobenzene	ND	10	"						
Anthracene	ND	10	"						
N-Nitrosodi-n-propylamine	ND	5.0	"						
1,2,4-Trichlorobenzene	ND	5.0	"						
1-Methylnaphthalene	ND	10	"						
2-Methylnaphthalene	ND	20	"						
4-Chloro-3-methylphenol	ND	10	"						
Acenaphthene	ND	10	"						
Benzo (a) anthracene	ND	10	"						
4-Nitrophenol	ND	10	"						
Benzo (b) fluoranthene	ND	10	"						
Benzo (k) fluoranthene	ND	10	"						
2,4-Dinitrotoluene	ND	10	"						
Pentachlorophenol	ND	10	"						
Benzo (g,h,i) perylene	ND	20	"						
Benzo (a) pyrene	ND	10	"						
Pyrene	ND	10	"						
Benzyl alcohol	ND	50	"						
Bis(2-chloroethoxy)methane	ND	10	"						
Bis(2-chloroethyl)ether	ND	5.0	"						
Bis(2-chloroisopropyl)ether	ND	20	"						
Bis(2-ethylhexyl)phthalate	ND	10	"						
4-Bromophenyl phenyl ether	ND	5.0	"						
Butyl benzyl phthalate	ND	10	"						
4-Chloroaniline	ND	20	"						
2-Chloronaphthalene	ND	10	"						
4-Chlorophenyl phenyl ether	ND	20	"						
Chrysene	ND	10	"						
Dibenz (a,h) anthracene	ND	10	"						
Dibenzofuran	ND	20	"						
Di-n-butyl phthalate	ND	5.0	"						

SunStar Laboratories, Inc.

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Benicia CA, 94510 Project Manager: Jim Gribi 10/08/14 17:03

Semivolatle Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 4100222 - EPA 3510C GCMS/ECD

Blank (4100222-BLK1)		Prepared: 10/02/14 Analyzed: 10/03/14	
1,2-Dichlorobenzene	ND	5.0	ug/l
1,3-Dichlorobenzene	ND	5.0	"
2,4-Dichlorophenol	ND	10	"
Diethyl phthalate	ND	10	"
2,4-Dimethylphenol	ND	5.0	"
Dimethyl phthalate	ND	10	"
4,6-Dinitro-2-methylphenol	ND	5.0	"
2,4-Dinitrophenol	ND	10	"
2,6-Dinitrotoluene	ND	20	"
Di-n-octyl phthalate	ND	10	"
Fluoranthene	ND	5.0	"
Fluorene	ND	10	"
Hexachlorobenzene	ND	20	"
Hexachlorobutadiene	ND	10	"
Hexachlorocyclopentadiene	ND	20	"
Hexachloroethane	ND	5.0	"
Indeno (1,2,3-cd) pyrene	ND	10	"
Isophorone	ND	10	"
2-Methylphenol	ND	10	"
4-Methylphenol	ND	20	"
Naphthalene	ND	5.0	"
2-Nitroaniline	ND	10	"
3-Nitroaniline	ND	10	"
4-Nitroaniline	ND	20	"
Nitrobenzene	ND	20	"
2-Nitrophenol	ND	10	"
N-Nitrosodiphenylamine	ND	10	"
N-Nitrosodimethylamine	ND	25	"
Phenanthrene	ND	10	"
2,4,5-Trichlorophenol	ND	20	"
2,4,6-Trichlorophenol	ND	10	"
2,3,4,6-Tetrachlorophenol	ND	10	"
2,3,5,6-Tetrachlorophenol	ND	10	"
1,4-Dinitrobenzene	ND	10	"
Pyridine	ND	10	"

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1090 Adam Street, Suite K Project Number: [none] Reported:
Benicia CA, 94510 Project Manager: Jim Gribi 10/08/14 17:03

Semivolatle Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 4100222 - EPA 3510C GCMS/ECD

Blank (4100222-BLK1)		Prepared: 10/02/14 Analyzed: 10/03/14						
Surrogate: 2-Fluorophenol	45.1	ug/l	100	45.1	9.97-110			
Surrogate: Phenol-d6	48.4	"	100	48.4	8.4-110			
Surrogate: Nitrobenzene-d5	65.5	"	100	65.5	14.7-110			
Surrogate: 2-Fluorobiphenyl	47.2	"	100	47.2	33.3-110			
Surrogate: 2,4,6-Tribromophenol	47.6	"	100	47.6	12.9-110			
Surrogate: Terphenyl-d14	46.3	"	100	46.3	15.8-136			
LCS (4100222-BS1)				Prepared: 10/02/14 Analyzed: 10/03/14				
Phenol	61.2	10	ug/l	100	61.2	12-89		
2-Chlorophenol	59.6	10	"	100	59.6	27-123		
1,4-Dichlorobenzene	49.7	10	"	100	49.7	36-97		
N-Nitrosodi-n-propylamine	65.8	5.0	"	100	65.8	41-116		
1,2,4-Trichlorobenzene	49.8	5.0	"	100	49.8	39-98		
4-Chloro-3-methylphenol	47.9	10	"	100	47.9	23-97		
Acenaphthene	63.4	10	"	100	63.4	46-118		
4-Nitrophenol	58.2	10	"	100	58.2	10-80		
2,4-Dinitrotoluene	63.0	10	"	100	63.0	24-96		
Pentachlorophenol	38.5	10	"	100	38.5	9-103		
Pyrene	58.6	10	"	100	58.6	26-127		
Surrogate: 2-Fluorophenol	55.5	"	100	55.5	9.97-110			
Surrogate: Phenol-d6	57.1	"	100	57.1	8.4-110			
Surrogate: Nitrobenzene-d5	51.3	"	100	51.3	14.7-110			
Surrogate: 2-Fluorobiphenyl	41.9	"	100	41.9	33.3-110			
Surrogate: 2,4,6-Tribromophenol	52.4	"	100	52.4	12.9-110			
Surrogate: Terphenyl-d14	41.2	"	100	41.2	15.8-136			
LCS Dup (4100222-BS1)				Prepared: 10/02/14 Analyzed: 10/03/14				
Phenol	59.5	10	ug/l	100	59.5	12-89	2.75	42
2-Chlorophenol	60.4	10	"	100	60.4	27-123	1.43	40
1,4-Dichlorobenzene	43.0	10	"	100	43.0	36-97	14.3	28
N-Nitrosodi-n-propylamine	61.2	5.0	"	100	61.2	41-116	7.15	38
1,2,4-Trichlorobenzene	45.3	5.0	"	100	45.3	39-98	9.50	28
4-Chloro-3-methylphenol	50.4	10	"	100	50.4	23-97	5.12	42
Acenaphthene	64.0	10	"	100	64.0	46-118	1.01	31
4-Nitrophenol	63.4	10	"	100	63.4	10-80	8.62	50
2,4-Dinitrotoluene	57.1	10	"	100	57.1	24-96	9.76	38

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Katherine RunningCrane

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Gribi Associates 1090 Adam Street, Suite K Benicia CA, 94510	Project: Maz Glass Project Number: [none] Project Manager: Jim Gribi	Reported: 10/08/14 17:03
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Semivolatle Organic Compounds by EPA Method 8270C - Quality Control

SunStar Laboratories, Inc.

Analyte	Result	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit	Notes
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Batch 4100222 - EPA 3510C GCMS/ECD

LCS Dup (4100222-bsd1)

Prepared: 10/02/14 Analyzed: 10/03/14

Pentachlorophenol	38.2	10	ug/l	100	38.2	9-103	0.887	50		
Pyrene	56.7	10	"	100	56.7	26-127	3.26	31		
Surrogate: 2-Fluorophenol	52.0		"	100	52.0	9.97-110				
Surrogate: Phenol-d6	54.5		"	100	54.5	8.4-110				
Surrogate: Nitrobenzene-d5	50.7		"	100	50.7	14.7-110				
Surrogate: 2-Fluorobiphenyl	55.2		"	100	55.2	33.3-110				
Surrogate: 2,4,6-Tribromophenol	48.8		"	100	48.8	12.9-110				
Surrogate: Terphenyl-d14	44.3		"	100	44.3	15.8-136				

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Gribi Associates 1090 Adam Street, Suite K Benicia CA, 94510	Project: Maz Glass Project Number: [none] Project Manager: Jim Gribi	Reported: 10/08/14 17:03
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Notes and Definitions

DET Analyte DETECTED
 ND Analyte NOT DETECTED at or above the reporting limit
 NR Not Reported
 dry Sample results reported on a dry weight basis
 RPD Relative Percent Difference

SunStar Laboratories, Inc.

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