

PROGRESS REPORT Quanta Resources, Lodi Street City of Syracuse, Onondaga County, New York DEC Site No. 7-34-013 / Index No. D7-00001-07-07 Plumley Project No. 2015127

January 2019

### **INTRODUCTION**

This report summarizes the remedial activities at the Quanta Resources – Syracuse Site since the last progress report (January 2017). Free product recovery efforts have been ongoing since March 2017, with the installation of absorbent socks in four monitoring wells with a history of free product accumulation. The Periodic Review Report (PRR) was submitted to the New York State Department of Environmental Conservation (DEC) in July 2018. The third post-remedial groundwater sampling event of selected wells was completed in December 2018.

### FREE PRODUCT RECOVERY

Absorbents have been maintained in the four wells with a history of free product accumulation (MW-1S, MW-2, MW-7 and MW-10) since March 2017. Quarterly inspections have been completed to assess the amount of oil absorbed in each well. During this period, a range of approximately 0.3 to 4.75 feet of free product accumulation was observed in well MW-1S and an estimated 0.98 gallons was recovered with the absorbents and bailers. Approximately <0.1 to 0.3 feet of free product accumulation well MW-7 and an estimated 0.21 gallons was recovered with the absorbents. No free product was observed or recovered from wells MW-2 and MW-10. Refer to the attached *Table 1A – Monitoring Well Groundwater Elevation* 

*Summary, Table 1B – Free Product Thickness* and *Table 1C – Monthly Free Product Recovery* for additional information.

#### **GROUNDWATER SAMPLING**

The depth to groundwater and free product thickness were measured in all of the wells and a groundwater sampling event was completed on December 5, 2018. Well MW-1S had a measurable accumulation of free product and was not sampled. Plumley personnel purged the wells and collected samples from wells MW-1D, MW-2, MW-5, MW-6, MW-7, MW-9, MW-10 and MW-12. Only wells MW-2 and MW-10 had odors after baling. No sheens or free product were observed in any of the wells sampled. Refer to the attached *Figure 1 – Site Plan* for sampling locations and *Groundwater Sampling Logs* for field observations.

Samples were submitted to SGS Accutest (SGS) for laboratory analysis of volatile organic compounds (VOCs) per United States Environmental Protection Agency (EPA) Method 8260 and polychlorinated biphenyls (PCBs) per EPA Method 8082. At the request of the New York State Department of Environmental Conservation (DEC), key monitoring well samples were also submitted to SGS for "emerging contaminants," including 1,4-Dioxane per EPA Method 8270SIM and Per- and Poly-Fluorinated Alkyl Substances (PFAS) Target Analyte List (TAL) per EPA Method 537M by ID. The results are discussed below.

#### **Groundwater Elevations**

The groundwater contours generated from elevation data collected on December 5, 2018 are shown on the attached *Figure 1 – Site Plan*. The elevation data indicate predominant groundwater flow directions generally to the west and southwest, similar to the last sampling

event in October 2017. Refer to the attached *Table 1A – Monitoring Well Groundwater Elevation Summary* for additional information.

#### **Analytical Results**

Analytical results showed total VOC concentrations ranging from 11 to 36 micrograms per liter ( $\mu$ g/L) in the wells tested. VOC concentrations were generally consistent with the results of the October 2017, December 2016 and September 2015 sampling events.

PCBs were detected in groundwater samples from well MW-2 and MW-10. Aroclors 1248, 1254 and 1260 were present at a total concentration of 3.05  $\mu$ g/L in MW-2. Aroclor 1254 was present at a concentration of 0.28  $\mu$ g/L in MW-10. These total concentrations were above the State groundwater standard<sup>1</sup> of 0.09  $\mu$ g/L. Well MW-2 had slight sheen during purging of the well and only an odor when sampled. MW-10 had an oil film present during purging of the well and an odor when sampled. These findings continue to be similar to the prior post-remedial sampling events.

PFAS were detected in groundwater samples from key monitoring wells MW-9 and MW-12. The PFAS reported in these two wells were at concentrations ranging from 1.31 to 23.6 nanograms per liter (ng/L). No PFAS were detected above the laboratory's method of detection limit in the sample collected from MW-10. No State standards or guidelines have been established for PFAS. The EPA has issued a health advisory of 70 ng/L, based on lifetime exposure to PFAS.

1,4-Dioxane was also tested in groundwater samples from key monitoring wells MW-9, MW-10 and MW-12. 1,4-Dioxane concentrations were reported at 2.34, 0.225 and 0.239  $\mu$ g/L,

<sup>&</sup>lt;sup>1</sup>DEC Division of Water Technical and Operational Guidance Series (TOGS) 1.1.1, *Ambient Water Quality Standards and Guidance Values and Groundwater Effluent Limitations*, dated June 1998 and Addenda.

respectively. There are no State or Federal standards or guidelines for 1,4-Dioxane in groundwater. The EPA has set a screening level<sup>2</sup> of 0.46  $\mu$ g/L in tap water.

Refer to the attached Figure 2 – VOC & PCB Groundwater Data Plan, Table 2 – Summary of Groundwater Analytical Results – VOCs and PCBs, Table 3 – Summary of Historical Groundwater Analytical Results – VOCs [Detections Only], Table 4 – Summary of Historical Groundwater Analytical Results – Total PCBs, Table 5 – PFAS and 1,4-Dioxane in Groundwater, Well Inspection Logs, Groundwater Sampling Field Logs and Laboratory Report for additional information.

#### CONCLUSIONS

We offer the following conclusions based on the groundwater sampling results:

- The VOC impacts to groundwater have not changed significantly from the 2017, 2016 and 2015 post-remedial sampling events.
- PCBs were present in groundwater above standards in only two onsite wells (MW-2 and MW-10) with a history of free product accumulation, similar to prior results.
- No significant migration of VOCS or PCBs to offsite wells is indicated.
- Free product accumulation was present in well MW-1S is consistent with past findings. The increase in accumulation observed in December 2018 was likely a result of wet weather in the fall of 2018. An estimated 0.98 gallons of free product were recovered

<sup>&</sup>lt;sup>2</sup>Technical Fact Sheet, 1,4-Dioxane; USEPA; November 2017.

from well MW-1S during the absorbent monitoring period, or approximately an average of 0.2 gallons per month. Absorbents in well MW-7 recovered 0.21 gallons during the period.

- Well RW-3 had trace amounts of free product in February, March and December of 2018. Measurements indicated a thickness of <0.1 feet. Well RW-3 has a history of sporadic free product impacts.</li>
- Although oil migration rates are low, continued accumulation of oil in some of the wells can be expected. Such findings are not necessarily indicative of a thick free product layer, but rather are more likely associated with the viscous oil gradually accumulating in the permeable well boring from an elevation above the water table.
- Three key wells were sampled for PFAS. Two of the three wells had detections of PFAS at concentrations ranging from 1.31 to 23.6 ng/L, which are well below the EPA health advisory concentration of 70 ng/L.
- All three key monitoring wells tested for 1,4-Dioxane had reportable concentrations of the compound. The reported concentrations were 2.34, 0.225 and 0.239 µg/L for wells MW-9, MW-10 and MW-12, respectively.

These results will be incorporated into the next PRR, which is due in July 2019.

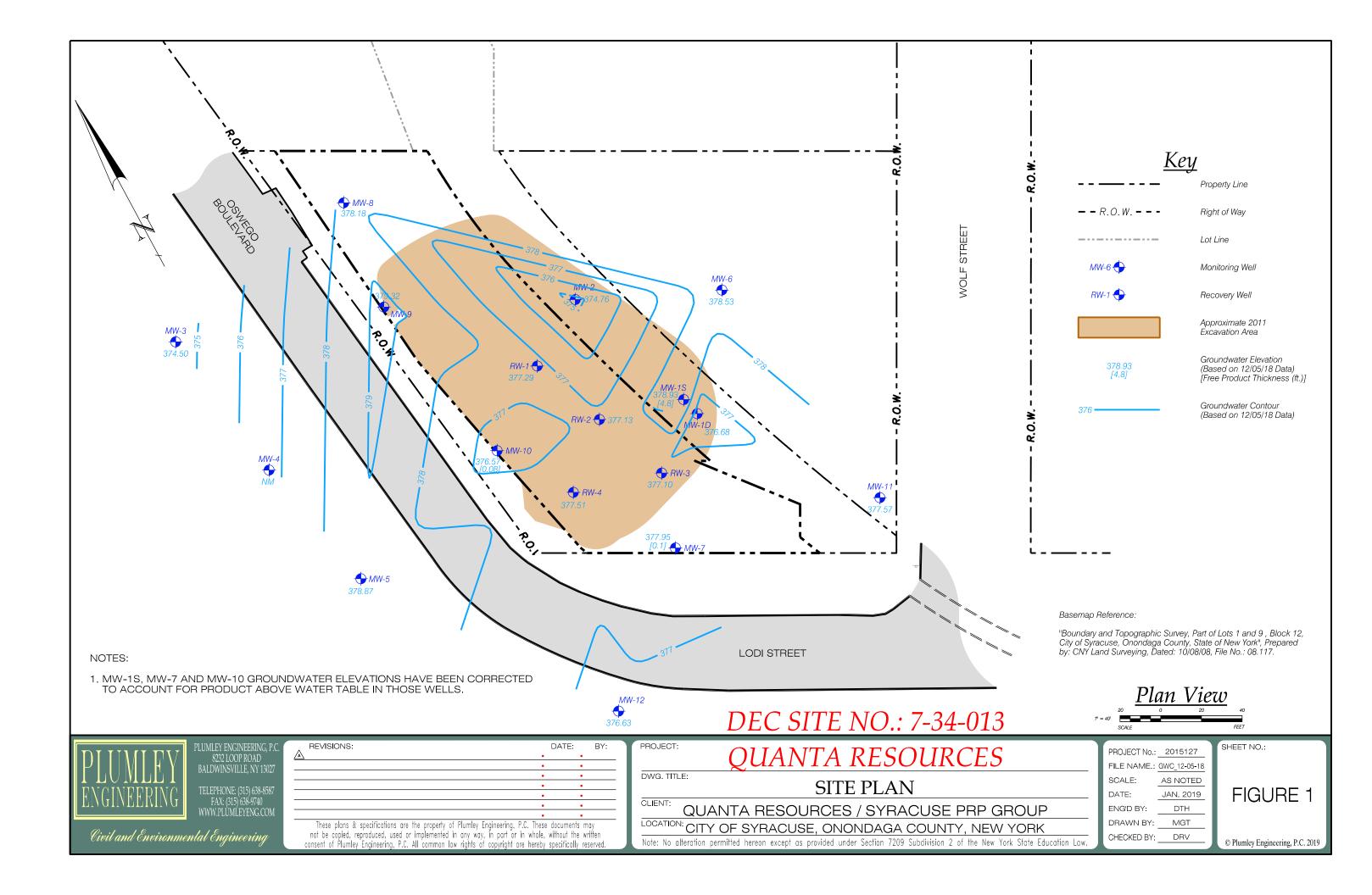
#### RECOMMENDATIONS

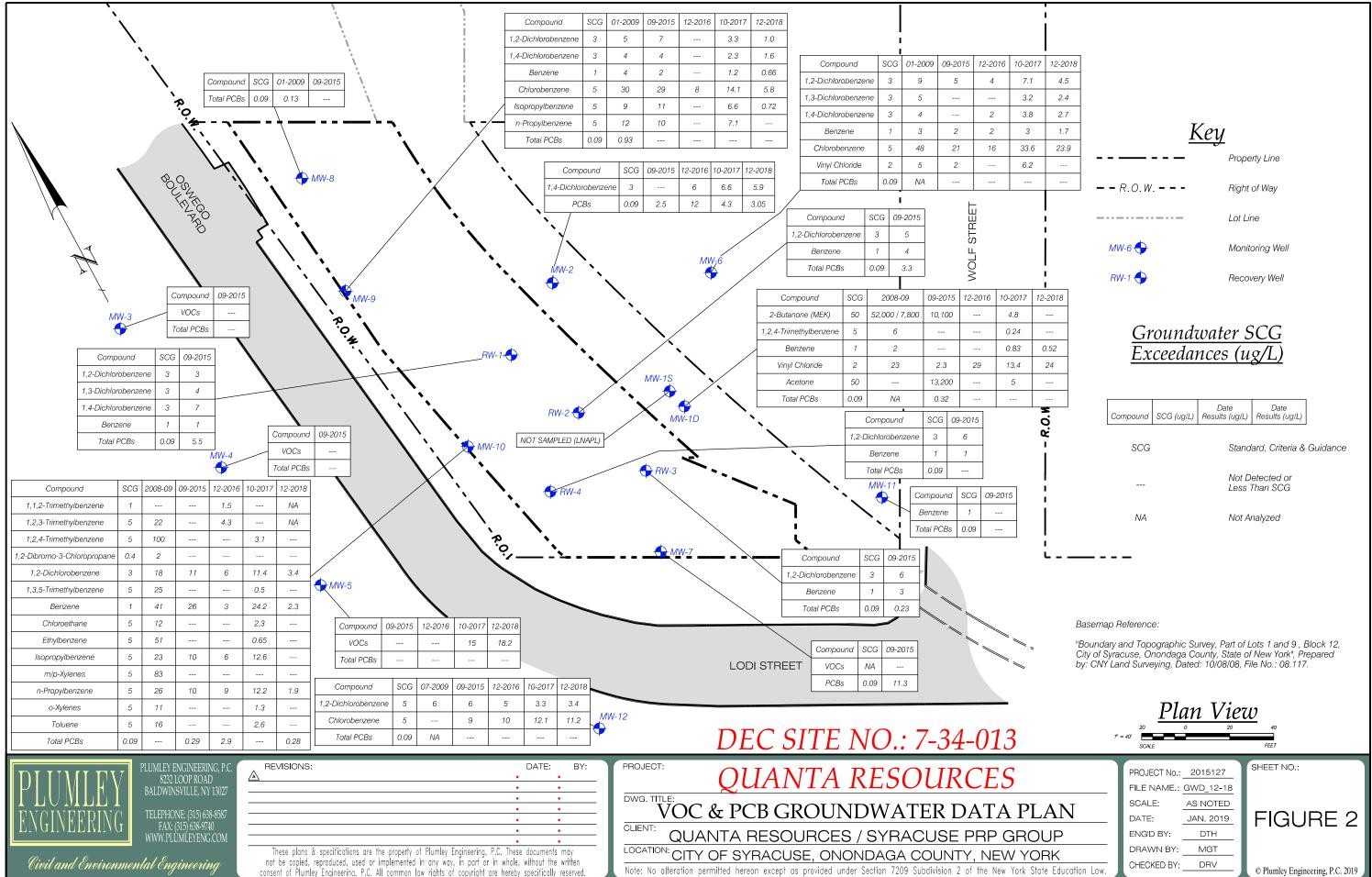
We recommend maintaining absorbent socks in the four wells with a history of free product accumulation (MW-1S, MW-2, MW-7 and MW-10) and well RW-3. Monitoring of these wells

for free product thickness and replacement of spent absorbents should be conducted quarterly. Other wells should be checked annually for free product.

We recommend continuing the groundwater sampling program but extending the time between sampling events to two years, given the consistency of the post-remedial results. The next sampling event would be in the summer of 2020 and would include wells MW-1S, MW-2, MW-5, MW-6, MW-7, MW-9, MW-10 and MW-12. Groundwater samples collected from these wells will be submitted for laboratory analysis of VOCs per EPA Method 8260 Site Specific Target Compound List (TCL) and PCBs per EPA Method 8082. No additional sampling and analysis of emerging contaminants appears to be warranted.

# FIGURES





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# **TABLES**

#### **TABLE 1A - MONITORING WELL GROUNDWATER ELEVATION SUMMARY**

WELL ID	MW-lD	MW-IS	MW-2	MW-3	MW-4	MW-5	MW-6	<b>MW-7</b>	MW-8	MW-9	<b>MW-10</b>	MW-11	MW-12	RW-1	<b>RW-2</b>	RW-3	RW-4
RISER ELEVATION	407.02	407.19	406.92	399.9	399.9	399.45	408.5	404.94	406.06	406.9	406.86	406.74	403.44				
GROUND ELEVATION	405.04	404.64	405.45	398.42	398.09	398.11	406.01	402.52	403.61	404.38	404	404.22	401.01				
<b>RISER ELEVATION post excavation</b> (1/1/2011)	407.23	404.56	405.36	399.9	399.9	401.12	408.5	402.08	404.59	406.91	403.61	406.74	403.43	404.61	404.08	403.5	402.92
GROUND ELEVATION post excavation (1/1/2011)	404.66	404.73	405.56	398.42	398.09	396.96	406.01	402.52	402.78	404.42	403.9	404.22	401.54	404.84	404.38	404.04	403.41
ELEVATIONS OF (Top)	365.04	370.64	377.45	373.42	366.09	376.11	387.51	389.02	386.11	386.88	384.5	384.72	381.51	381.88	381.54	384.04	383.41
SCREEN INTERVAL (Bottom)	360.04	365.64	367.45	358.42	356.09	361.61	367.51	373.02	372.11	370.88	369.5	369.72	366.51	361.88	361.54	364.04	363.41
BOTTOM OF BORING ELEVATION	357.04	365.64	367.45	356.92	355.59	359.11	367.01	372.52	371.61	370.38	369.00	369.22	366.01	361.88	361.54	364.04	363.41
DATE INSTALLED	11/18/91	11/25/91	11/21/91	11/26/91	11/25/91	11/27/91	12/18/08	12/11/08	12/09/08	12/10/08	12/16/08	06/25/09	07/09/09	12/16/08	06/25/09	01/02/10	07/12/10
DIAMETER (Inches)	2	2	2	2	2	2	2	2	2	2	2	2	2	2	2	2	2
CASING MATERIAL	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC
SCREEN MATERIAL	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC	PVC
SLOT SIZE (Inches)	0.010	0.010	0.010	0.010	0.010	0.010	0.020	0.020	0.020	0.020	0.020	0.020	0.020	0.020	0.020	0.020	0.020
DATE*	01010	01010	01010	0.010	0.010	0.010	0.020	GROUNDV				0.020	0.020				
02/06/1992	374.45	376.81	377.8	374.03	374.00	378.46	NI	NI	NI	NI	NI	NI	NI	NI	NI	NI	NI
04/15/1992	375.37	377.77	378.62	374.96	374.89	378.56	NI	NI	NI	NI	NI	NI	NI	NI	NI	NI	NI
03/10/2008**	374.37		376.58 (2.3')	373.51	373.29	377.33	NI	NI	NI	NI	NI	NI	NI	NI	NI	NI	NI
03/12/2008*	374.5	NM	NM	373.43	373.33	377.06	NI	NI	NI	NI	NI	NI	NI	NI	NI	NI	NI
12/16/2008	NM	NM	NM	NM	NM	NM	NI	375.36	377.56	377.59	NI	NI	NI	NI	NI	NI	NI
12/18/2008	NM	NM	NM	NM	NM	NM	NI	375.61 (.04')	378.05	377.55	377 0	NI	NI	NI	NI	NI	NI
12/23/2008	NM	NM	NM	NM	NM	NM	377.05	375.60 (.16')	377.73	377.53	376.8 0	NI	NI	NI	NI	NI	NI
01/05/2009*	375.58	NM	NM	374.6	374.55	376.53	377.41	376.41 (.26')	378.3	378.26	377.5 0	NI	NI	NI	NI	NI	NI
01/23/2009**	374.41		376.78 (2.07')	374.14	374.01	375.65	375.77	375.22 (.44')	377.5	376.99	376.5 0	NI	NI	NI	NI	NI	NI
06/25/2009**	374.37		375.94 (1.86')	373.79	373.69	375.7	375.41	375.06 (.38')	377.64	376.67	376.32 (.25')	NI	NI	NI	NI	NI	NI
06/29/2009	374.36		376.10 (1.51')	373.72	373.66	375.97	375.22	374.86 (.64')	377.37	376.61	376.15 (.29')	NI	NI	NI	NI	NI	NI
07/14/2009**	374.16	374.87 (2.04')	375.81	373.61	373.54	375.44	374.99	371.59	376.97	376.25	375.6 (.85')	NI	373.94	NI	NI	NI	NI
09/22/2015**	378.83	379.91 (3.0')	375.56	372.7	373.9	380.12	380.3	373.68	378.39	381.51	375.81	377.34	377.83	376.11	381.38	374.9	379.12
12/08/2016**	375.48	377.36 (3.0')		374.51	NM	377.86	377.07	377.00 (.38')	378.21	378.08	375.94 (.08')	376.22	375.67	378.03	377.81	377.14	377.65
12/16/2016*	375.25	377.17 (3.0')	377.68	NM	375.2	377.87	376.54	376.79	NM	377.59	375.03	NM	375.20	NM	NM	NM	NM
03/01/2017	NM	378.06 (4.0')		NM	NM	NM	NM	378.11 (.27')	NM	NM	378.99 (.09')	NM	NM	NM	NM	NM	NM
04/25/2017 05/26/2017	NM NM	377.97 (2.5')	378.26	NM NM	NM	NM	NM	377.48 (1')	NM	NM NM	378.41	NM NM	NM NM	NM NM	NM NM	NM	NM NM
06/29/2017	375.47	376.64 (1.3') 375.95 (0.7')	377.71 377.95	374.2	NM NM	NM 377.15	NM 376.51	377.21 (.15') 376.49	NM 378.02	377.8	377.82 375.81	375.93	NM 375.52	377.92	377.04	NM 376.91	377.21
07/29/2017	3/3.4/ NM	376.05 (0.7)	377.95	374.2 NM	NM	NM	376.31 NM	376.49	378.02 NM	377.8 NM	375.81	NM	NM	378.44	377.96	376.91	377.21
09/08/2017	NM	378.52 (0.8')	378.43	NM	NM	NM	NM	377.92 (<0.1')	NM	NM	378.87	NM	NM	378.21	379.4	377.12	377.6
10/05/2017	374.19	376.64 (1.5')	377.71	372.15	NM	377.35	374.68	377.21 0.2	376.19	375.92	378.87	374.69	374.50	376.6	377.42	375.43 0.2	375.59
02/28/2018	574.19 NM	375.95 (1.8')	377.95	NM	NM	401.12	574.08 NM	376.49 0.3	NM	NM	375.81	NM	NM	376.61	377.42	375.52 0	375.77
05/30/2018	375.28	376.05 (0.8')	378.45	373.99	NM	376.91	376.38	377.07 0.1	377.88	377.6	376.39	375.75	375.39	377.72	376.75	375.32 0 376.78 0	377.05
09/28/2018	NM	378.52 (0.3')	377.37	NM	NM	401.12	NM	377.92 0	377.88	NM	378.87	NM	NM	378.53		377.49 0	377.75
12/05/2018	376.68	378.93 (4.8')	374.76	374.5	NM	378.87		377.95 0.1	378.18	379.32	376.57	377.57	376.63	377.29		377.10 0	377.51
	570.00	570.75 (0.7)	517.10	577.5	LVIVI	570.07	570.55	511.75 0.1	570.10	517.54	570.57	511.51	570.05	511.47	511.15	577.10 0	511.51

Notes:

(1.1') indicates measured free product thickness in feet.

All wells were re-surveyed on 01/05/09 by Plumley Engineering and those elevations were used for all groundwater data from 03/10/08 to 2012. Wells re-surveyed after excavation, those elevations were used for all groundwater data from 2012 to present.

\*\*Wells contained free product layers on the water column. A Corrected Depth To Water (CDTW) calculation was used to estimate the groundwater level without the free product using this equation: CDTW = Static DTW - (PxG); where P = Measured Product thickness (which is notated in parenthesis) and G = Specific Gravity. Specific Gravity is currently estimated to be 0.85 based on field observations and published values.

\*Groundwater sampling date.

NI Not installed

NM Not measured

All elevations reported in feet above mean sea level.

#### TABLE 1B - FREE PRODUCT THICKNESS (FEET)

Date	MW-1S	MW-2	MW-7	MW-10	RW-1	RW-2	RW-3	RW-4	MW-1D	MW-3	MW-4	MW-5	MW-6	MW-8	MW-9	MW-11	MW-12
07/14/09	2.04	1.80	0.90	0.85	NI	NI	NI	NI									
2011	2.0.1	100	0.50	0.00					eted Remedial Ex								
09/20/12	3.23	1.07	4.03	2.09		0.05	0.76										
09/20/12				,					System Startup		1	<b>I</b>		1	<b>I</b>		
09/27/12	3.20	1.51	3.21	1.68		0.14	0.19										
10/04/12	4.26	1.39	4.85	2.05		0.09	0.27										
10/12/12	4.25	2.21	4.49	1.69		0.66	0.95										
11/15/12	NA	0.77	NA	1.5		NA	NA										NM
12/28/12	6.21	1.01	2.92	1.32		0.31	NA										
01/30/13	6.4	0.29	0.33	0.87		0.32	0.13										
02/22/13	4.76	0.13	2.01	1.37		0.18	0.19										
03/28/13	3.41	0.13	2.31	1.37		0.68	0.3										
04/30/13	1.14	0.06	1.40	0.96													
05/30/13	1.62	0.77	1.36	0.95		NA	0.21										
06/21/13	2.29	0.13	0.82	0.91		0.55	0.43										
07/17/13	1.70	0.09	1.56	0.53		< 0.01	0.21										
08/15/13	0.50	0.20	0.11	0.30		0.02	0.02	0.02									
09/25/13	3.00	0.05	0.50	0.75		0.01	0.01										
10/30/13	3.00	0.01	0.50	0.75		0.05	NA										
11/21/13	3.00	0.08	1.00	0.33			0.01										
12/31/13	0.60	0.10	0.10	0.20			0.01	0.01									
01/31/14	3.00	NM	NM	NM		NM	NM	0.01									
01/31/14							ommence Absorb	ent Oil Recover	y Program (Free	Product Thickne	ss Not Measural	ole)			<u>.</u>	<u>.</u>	
02/02/14	0.00	0.00	0.00	0.00		0.00		0.00									
02/03/14	0.00	0.00	0.00	0.00		0.00	0.00	0.00									
02/04/14	0.00	0.00	0.00	0.00		0.00	0.00	0.00									
02/05/14	0.00	0.00	0.00	0.00		0.00	0.00	0.00									
09/22/15								А	bsorbents Remov	red					-	-	
09/22/15	3.00																
12/08/16	2.70	0.08	0.38	0.08													
02/16/16	3.50	0.25															
04/02/16	3.75	0.20															
07/12/16	4.33	0.25															
10/14/16	4.75	0.50	0.10	0.10													
12/08/16	3.66	0.10	0.40	0.10													
12/14/16	2.70	0.08	0.38	0.08													
03/01/17									Wells MW1S, M	,							
03/01/17	4.00	0.02	0.27	0.09	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
04/25/17	2.50	< 0.1	1.00	< 0.1	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
05/26/17	1.30	< 0.1	0.15	< 0.1	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
06/29/17	0.70	< 0.1	0.20	< 0.1							NM						
07/29/17	0.30	< 0.1	< 0.1	< 0.1					NM	NM	NM	NM	NM	NM	NM	NM	NM
09/08/17	0.80	< 0.1	0.10	< 0.1					NM	NM	NM	NM	NM	NM	NM	NM	NM
10/05/17	1.50	< 0.1	0.20	< 0.1			0.20	NM									
02/28/18	1.75	< 0.1	0.30	<0.2			< 0.1	NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
05/30/19	0.80	< 0.1	< 0.1	< 0.1			< 0.1	NM									
09/28/18	0.30							NM	NM	NM	NM	NM	NM	NM	NM	NM	NM
12/05/18	4.75	< 0.1	0.10	< 0.1			<0.1	NM									

Notes:

--- Not Present

NI Well not installed

NA Oil-water interface probe malfunction

NM Not measured

Free product measurements taken with an oil-water interface probe.

Plumley Engineering, P.C.

#### TABLE 1C - MONTHLY FREE PRODUCT RECOVERY\* (GALLONS)

Date	MW-1S	MW-2	<b>MW-7</b>	<b>MW-10</b>	RW-1	RW-2	RW-3	RW-4	Total	Cumulative Total
Sep-12		Syst	em Startup	and Montl	nly Manua	l Bailing B	legun			
Oct-12	1.32	0.26	0.92	0.40					2.9	2.9
Nov-12	0.69	0.13	0.79	0.18					1.8	4.7
Dec-12	0.99	0.08	0.79	0.20					2.1	6.8
Jan-13	0.90	0.01	0.03	0.11					1.0	7.8
Feb-13	1.82	0.01	0.42	0.13					2.4	10.2
Mar-13	1.77		0.29	0.11					2.2	12.4
Apr-13	1.43		0.17	0.11					1.7	14.1
May-13	0.54	0.03	0.16	0.05					0.8	14.8
Jun-13	0.29		0.16	0.05					0.5	15.4
Jul-13	0.21		0.26						0.5	15.8
Aug-13	0.20	0.01	0.11	0.01					0.3	16.2
Sep-13	0.26	0.01	0.11	0.05					0.4	16.6
Oct-13	0.20	0.01	0.03	0.05					0.3	16.9
Nov-13	0.21	0.01	0.08	0.03					0.3	17.2
Dec-13	0.16	0.01	0.01	0.03					0.2	17.4
Jan-14	0.26								0.3	17.7
Jan-14			Placed	d Absorben	ts Into All	Wells	-			
Feb-14	1.01	0.13	0.26	0.26	0.08	0.30	0.12	0.08	2.24	19.9
Mar-14	0.26		0.13	0.13	0.06		0.05		0.63	20.6
Apr-14			0.07	0.10	0.01	0.08	0.02		0.28	20.8
May-14			0.05	0.01	0.01	0.03	0.01	0.01	0.13	21.0
Jun-14	0.02	0.03	0.02	0.07	0.000	0.03	0.08	0.01	0.25	21.2
Jul-14	0.14	0.04	0.02	0.05	0.00	0.03	0.05	0.01	0.34	21.6
Aug-14	0.20	0.08	0.01	0.13	0.00	0.04	0.05	0.00	0.51	22.1
Sep-14	0.27	0.01	0.03	0.13	0.01	0.04	0.03	0.02	0.54	22.6
Oct-14	0.27	0.03	0.00	0.13	0.00	0.09	0.02		0.54	23.1
Nov-14	0.27	0.08	0.00	0.13	0.19	0.19			0.86	24.0
Dec-14	0.27	0.03	0.05	0.09	0.08	0.02			0.54	24.5
Jan-15	0.27	0.00	0.07	0.09	0.11	0.11	0.04	0.00	0.68	25.2
Mar-15	0.27	0.00	0.01	0.01	0.08	0.08	0.11	0.02	0.57	25.8
Apr-15	0.08	0.07	0.03	0.01	0.04	0.11	0.09	0.17	0.60	26.4
May-15	0.07	0.03	0.05	0.04	0.02	0.11	0.04	0.00	0.36	26.7
Jun-15	0.12	0.00	0.03	0.04	0.05	0.08	0.04	0.00	0.35	27.1
Aug-15	0.02	0.03	0.02	0.02	0.02	0.11	0.11	0.00	0.34	27.4
Mar-17	А	borbents re	einstalled i	n wells MV	V-1S, MW	'-2, MW-7	and MW-1	0		
Apr-17	0.27								0.27	27.7
May-17	0.27				· · · · · ·				0.27	28.0
Jun-17	0.27		0.20		· · · · · ·				0.47	28.4
Jul-17	0.27				· · · · · ·				0.27	28.7
Sep-17	0.40								0.40	29.1
Oct-17	0.40		0.20						0.60	29.7
Feb-18	0.27		0.07						0.34	30.1
May-18	0.03		0.01						0.04	30.1
Sep-18	0.03								0.03	30.1
Dec-18	0.65		0.13						0.78	30.9
Total	17.0	1.1	5.7	3.0	0.8	1.4	0.9	0.3	30.1	l

#### Notes:

\*Based on estimate in each bailer during bailing program and spent absorbent length during absorbent program. Blank indicates not present/removed. For wells not listed, free product is not present.

#### **TABLE 2 - SUMMARY OF GROUNDWATER ANALYTICAL RESULTS - VOCs and PCBs**

**Date Sampled:** December 5, 2018

Lab Sample ID:         Nature 179316-9         JC79316-9         JC7916-9         JC7916-9         JC7916-9	Client Sample ID:	Unit	State	MW-1-D	MW-2	MW-5	MW-6	MW-9	<b>MW-10</b>	MW-12
Actions         ug/L         -1         ND (6.0)         ND (6.	Lab Sample ID:	Umt	<b>Standard</b> <sup>1</sup>	JC79316-6	JC79316-5	JC79316-4	JC79316-3	JC79316-1	JC79316-8	JC79316-7
Binance         ygL         1         0.23         1.4         ND (A3)         1.7         0.66         2.3         0.04           Bronnobencem         ygL         5         ND (0.53)         ND (0.53)         ND (0.53)         ND (0.55)         ND (0.57)         ND (0.57) <td></td> <td></td> <td></td> <td>MS Vola</td> <td>tiles (SW84</td> <td>6 8260C)</td> <td>-</td> <td>-</td> <td>-</td> <td></td>				MS Vola	tiles (SW84	6 8260C)	-	-	-	
Bronnehursence         ight         5         ND (0.55)         ND (0.55)         ND (0.55)         ND (0.45)         ND (0.5)         ND (0.5) <thnd (0.5)<="" th="">         ND (0.5)         <thn< td=""><td>Acetone</td><td>µg/L</td><td>-</td><td>ND (6.0)</td><td>ND (6.0)</td><td>ND (6.0)</td><td>ND (6.0)</td><td>ND (6.0)</td><td>ND (6.0)</td><td>ND (6.0)</td></thn<></thnd>	Acetone	µg/L	-	ND (6.0)	ND (6.0)	ND (6.0)	ND (6.0)	ND (6.0)	ND (6.0)	ND (6.0)
Biomschlacementane         up1         S         ND (0.48)         ND (0.49)         ND (0.47)         ND (0.49)         ND (0.47)         ND	Benzene	µg/L	1	0.52	1.4	ND (0.43)	1.7	0.66	2.3	0.46 J
Bromoshame         μpT         -         ND 0.58         ND 0.58         ND 0.58         ND 0.58         ND 0.68         ND 0.65         ND 0	Bromobenzene	µg/L	5	ND (0.55)	ND (0.55)	ND (0.55)	ND (0.55)	ND (0.55)	ND (0.55)	ND (0.55)
Bromoscham         μgL         -         N D 0.63         N D 0.65         N D 0.63         N D	Bromochloromethane	µg/L	5	ND (0.48)	ND (0.48)	ND (0.48)	ND (0.48)	ND (0.48)	ND (0.48)	ND (0.48)
Brommerhume         μpT         5         NT (1.67         NT (1.67         NT (1.67         NT (1.67         NT (1.67)         NT (1.67) <td>Bromodichloromethane</td> <td>µg/L</td> <td>-</td> <td>ND (0.58)</td> <td>ND (0.58)</td> <td>0.82 J</td> <td>ND (0.58)</td> <td>ND (0.58)</td> <td>ND (0.58)</td> <td>ND (0.58)</td>	Bromodichloromethane	µg/L	-	ND (0.58)	ND (0.58)	0.82 J	ND (0.58)	ND (0.58)	ND (0.58)	ND (0.58)
2-Buamons (MTK)         µgT         -         ND (6.9)         ND (6.6)         ND (6.6) <td< td=""><td>Bromoform</td><td>µg/L</td><td>-</td><td>ND (0.63)</td><td>ND (0.63)</td><td>ND (0.63)</td><td>ND (0.63)</td><td>ND (0.63)</td><td>ND (0.63)</td><td>ND (0.63)</td></td<>	Bromoform	µg/L	-	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)	ND (0.63)
n-Baughtenzene         ngT.         5         ND (0.52)         ND (0.55)         ND (0.56)         ND (0.57)         ND (0	Bromomethane	µg/L	5	ND (1.6) <sup>a</sup>	ND (1.6) <sup>a</sup>	ND (1.6) <sup>a</sup>	ND (1.6) <sup>a</sup>	ND (1.6) <sup>a</sup>	ND (1.6) <sup>a</sup>	ND (1.6) <sup>a</sup>
ses-Enziphenzene         sgl.         S         ND (0.62)         ND (0.62)         ND (0.62)         ND (0.62)         ND (0.63)         ND (0.63)         ND (0.65)         ND (0.55)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.55)         ND	2-Butanone (MEK)	µg/L	-	ND (6.9)	ND (6.9)	ND (6.9)	ND (6.9)	ND (6.9)	ND (6.9)	ND (6.9)
intre-Burghbanzenc         ingL         5         ND (0.69)         ND (0.69)         ND (0.69)         ND (0.69)         ND (0.95)         ND (0.75)         ND (0.76)         N	n-Butylbenzene	µg/L	5	ND (0.52)	ND (0.52)	ND (0.52)	ND (0.52)	ND (0.52)	0.63 J	ND (0.52)
Carbon disulfide $\mu_{g}I_{L}$ 60         ND (0.95)         ND (0.95)         ND (0.95)         ND (0.95)         ND (0.55)         ND (0.56)	sec-Butylbenzene	µg/L	5	ND (0.62)	ND (0.62)	ND (0.62)	ND (0.62)	ND (0.62)	1.3 J	1.4 J
Carbon tetrachloride         jpg1         5         ND (0.55)	tert-Butylbenzene	µg/L	5	ND (0.69)	ND (0.69)	ND (0.69)	ND (0.69)	ND (0.69)	ND (0.69)	0.96 J
Carbon tetrachloride         μg/L         5         ND (0.53)         ND (0.55)         ND (0.56)         ND (0.57)         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.51)	Carbon disulfide		60	ND (0.95)	ND (0.95)	ND (0.95)	ND (0.95)	ND (0.95)	ND (0.95)	ND (0.95)
Chlorobacaca         μg/L         5         0.07.1         0.07.3         ND (0.53)         ND (0.73)         ND (0.75)	Carbon tetrachloride	1	5	ND (0.55)	ND (0.55)	ND (0.55)	ND (0.55)	ND (0.55)	ND (0.55)	ND (0.55)
Chloroschame         µgL         5         ND (0.73)         ND (0.74)         ND (0.74)	Chlorobenzene	1	5	0.57 J	0.97 J	ND (0.56)	23.9	5.8	ND (0.56)	11.2
Chloroform         µg/L         7         ND (0.50)         ND (0.51)         ND (0.50)         ND (0.51)         ND (0.50)         ND (0.50)         ND (0.51)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.51)         ND (0.51)         ND (0.52)         ND (0.52)         ND (0.52)         ND (0.52)         ND (0.54)         ND (0.54)         ND (0.55)         ND (0.54)         ND (0.55)         ND (0.57)         ND (0.51)         ND (0.55)         ND (0.55)<	Chloroethane		5	ND (0.73)			ND (0.73)	ND (0.73)	· · · · ·	ND (0.73)
Chloromethane         µg/L         5         ND (0.76)         ND (0.77)         ND (0.7	Chloroform		7		ND (0.50)		· /	× /	× /	· · · · · ·
o-Chlorotoluce $\mu g'L$ 5         ND (0.63)         ND (0.64)         ND (0.64)         ND (0.64)         ND (0.65)         ND (0.66)         ND (0.66)         ND (0.66)         ND (0.65)         ND (0.64)         ND (0.21)         ND (0.56)         ND (0.55)         ND (0.55)         ND (0.55)         ND (0.55)         ND (0.55)         ND (0.55)         ND (0.56)         ND (0.57)         ND (0.51)         Ad-5         I         Z.6         Ad-1           1.4 Dichlorobenzene $\mu g'L$ 5         ND (0.57)         S         ND (0.57)         ND (1.47)         ND (1.47)         ND (1.47)         ND (1.47)         ND (0.57)	Chloromethane		5	· · · ·	. ,		``´´	× /	<b>`</b>	· · · · ·
p-Chlorotoluce $\mu g'L$ 5         ND (0.60)         ND (0.61)         ND (0.71)         ND (0.71)         ND (0.72)         ND (0.72)         ND (0.72)         ND (0.73)         ND (0.73)         ND (0.73)         ND (0.74)         ND (0.75)         ND						· · · · · ·	· · · · ·	· · · · ·	× /	
12-Dibrome-3-chloropropune         µg/L         0.04         ND (1.2)         ND (1.4)         ND (0.56)         ND (0.56)         ND (0.50)         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.51)         ND (1.4)         ND		1		<b>`</b>	, , , , , , , , , , , , , , , , , , ,		<b>`</b>	· · · · ·	<b>`</b>	· · · · ·
Dihromechloromethane         μg/L         -         ND (0.56)         ND (0.57)         ND (0.51)         ND (0.51)         ND (0.51)         S.5         ND (0.51)         S.5         ND (0.51)         S.5         ND (0.51)         S.5         ND (0.51)         ND (0.51)         S.5         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.57)         ND (0.50)         ND (0.51)         ND	1			( )	× /		× /	~ /	· · · /	· · · · · ·
1.2-Dibromoethane         µg/L         0.0006         ND (0.48)         ND (0.41)         ND (0.51)         2.4         0.621         ND (0.48)         ND (0.48)         ND (0.47)         ND (1.47)         ND (0.50)         ND (0.51)         ND (0.51	* *			× /		· /			× /	
1.2-Dichlorobenzene         μg/L         3         0.59         3.1         ND (0.53)         4.5         1         2.6         3.4           1.3-Dichlorobenzene         μg/L         3         ND (0.54)         2.5         ND (0.54)         2.4         0.621         ND (0.54)         ND (0.51)         5.9         ND (0.51)         5.9         ND (0.51)         5.9         ND (0.51)         5.9         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.57)         ND (0.57)         ND (0.57)         ND (0.57)         ND (0.57)         ND (0.59)         ND (0.59)         ND (0.59)         ND (0.59)         ND (0.51)         ND (0.52)         ND (0.51)         ND (0.5				· /	. ,	. ,		× /	```´	
1.3-Dichlorobenzene $\mu g/L$ 3         ND (0.54)         2.5         ND (0.51)         2.7         ND (0.51)         ND (0.51)         LA         0.61         1.1         1.8           Dichlorodihuromethane $\mu g/L$ 5         ND (0.51)         S.9         ND (0.57)         ND (0.59)         ND (0.51)         ND (0.52)		-		× /	, <i>,</i> ,	· · · ·	. ,	1	× /	
1.4-Dichlorobenzene         µg/L         3         ND (0.51)         5.9         ND (0.51)         2.7         1.6         1.1         1.8           Dichlorodffluoromethane         µg/L         5         ND (1.4)         ND (0.57)         ND (0.57)         ND (0.57)         ND (0.57)         ND (0.50)         ND (0.51)         ND (0.52)         ND (0.43)								0.62 I		
Dicklorodifluoromethane         µg/L         5         ND (1.4)'         ND (0.5)'         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.51)         ND (0.52)         ND (0.51)         <	-									
1.1-Dichloroethane $\mu g'L$ 5         2.3         0.73 J         ND (0.57)         0.61 J         ND (0.57)         ND (0.59)         ND (0.59)         ND (0.59)         ND (0.59)         ND (0.59)         ND (0.51)         ND (0.52)         ND (0.51)         ND (0.51)<										
1.2-Dichloroethane $\mu g'L$ 0.6         ND (0.60)         ND (0.60)         ND (0.60)         ND (0.60)         ND (0.60)         ND (0.60)         ND (0.50)         ND (0.51)         ND (0.52)         ND (0.43)				· /	· · · · ·		, <i>, ,</i>	× /		· /
1.1-Dichloroethene         µg/L         5         ND (0.59)         ND (0.59)         ND (0.59)         ND (0.59)         ND (0.59)         ND (0.51)         ND (0.52)         ND						· · · ·		· · · · ·	```´	· · · ·
1,2-Dichloroethene (total)         µg/L         -         1.5         1         ND (0.51)         ND (0.52)         ND (0.51)         ND (0.47)         ND (0.43)         ND (0.43)         ND (0.43)         ND (0.43)         ND (0.47		1		· · · ·	<b>``</b>	× /	· /	× /	. ,	· · · ·
1,2-Dichloropropane         μg/L         1         ND (0.51)         ND (0.52)         ND (0.53)         ND (0.43)	,			× /	1	( )	· · · · ·	× /	× /	. ,
1.3-Dichloropropane $\mu g' L$ 5ND (0.43)ND (0.52)ND (0.43)ND (0.43) <td></td> <td></td> <td></td> <td></td> <td>ND(0.51)</td> <td>· /</td> <td>· /</td> <td>, , ,</td> <td></td> <td>· · · · ·</td>					ND(0.51)	· /	· /	, , ,		· · · · ·
2.2-Dichloropropane $\mu g/L$ 5         ND (0.52)         ND (0.53)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.43)         ND (0.43)         ND (0.43)         ND (0.43)         ND (0.56)         <		1		<b>`</b>	```´	. ,	. ,		``´´	· /
1.1-Dichloropropene $\mu g'L$ -         ND (0.82)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.43)         ND (0.50         ND (0.50         ND (0.56)         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.51) <th< td=""><td></td><td></td><td></td><td>× /</td><td>``´´</td><td>· · · ·</td><td>· · · ·</td><td>× /</td><td>``´´</td><td>``´´</td></th<>				× /	``´´	· · · ·	· · · ·	× /	``´´	``´´
cis-1,3-Dichloropropene         µg/L         -         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.47)         ND (0.43)         ND (0.45)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.56)         ND (0.50)         ND (0.55)         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.51)         ND (0.48)         ND (0.48)         <				· /	, , ,	· · · · · ·	· /	, <i>, ,</i>	``´´	
trans-1,3-Dichloropropene         μg/L         -         ND (0.43)         ND (0.50)         ND (0.50)         ND (0.50)         ND (0.56)         ND (0.56)         ND (0.56)         ND (0.56)         ND (0.56)         ND (0.51)         ND (0.43)         ND (0.43)         ND (0.51)		1		<b>`</b>	, , , , , , , , , , , , , , , , , , ,	. ,	<b>`</b>	· · · · ·	```´	. ,
Ethylbenzene $\mu g/L$ 5ND (0.60)ND (0.50)ND (0.51)ND				· /	``´´		· /	```´´	``´´	· /
Hexachlorobutadiene $\mu g/L$ 0.5ND (0.56)ND (0.57)ND (0.57)ND (0.57)ND (0.57)ND (0.57)ND (0.57) </td <td></td> <td></td> <td></td> <td>× /</td> <td>``´´</td> <td>· · · ·</td> <td>· · · ·</td> <td>× /</td> <td>```´</td> <td>. ,</td>				× /	``´´	· · · ·	· · · ·	× /	```´	. ,
Isopropylbenzeneµg/L5ND (0.65)ND (0.65)ND (0.65)ND (0.65)0.72 J2.33.9p-Isopropylolueneµg/L5ND (0.66)ND (0.66)ND (0.66)ND (0.66)ND (0.66)ND (0.66)ND (0.66)ND (0.66)ND (0.61)ND (0.51)ND (0.51) <td>·</td> <td></td> <td></td> <td>× /</td> <td>. ,</td> <td>· /</td> <td>· · · ·</td> <td>· · · · ·</td> <td><b>`</b></td> <td>· /</td>	·			× /	. ,	· /	· · · ·	· · · · ·	<b>`</b>	· /
p-Isopropyltolueneμg/L5ND (0.66)ND (0.51)ND (0.51) <th< td=""><td></td><td></td><td></td><td>× /</td><td>. ,</td><td>× /</td><td>· · · ·</td><td></td><td>× /</td><td></td></th<>				× /	. ,	× /	· · · ·		× /	
Methyl Tert Butyl Ether $\mu g/L$ 10ND (0.51)ND (0.51	1 10			<b>`</b>	<b>`</b>	· · · ·				
Methylene bromide         μg/L         5         ND (0.48)         ND (0.50)         ND (0.50)         ND (0.50         ND (0.50)         ND (0.50         ND (				· /	` <i>´</i>	. /	· · · /			· /
Methylene chloride         μg/L         5         ND (1.0)         ND (0.8)         ND (0.80         ND (0.80         ND (0.80         ND (0.60				× /	. ,		· · · ·	× /	``´´	. ,
Naphthalene         μg/L         -         ND (0.98)         ND (0.90)         ND (0.60)         ND (0.60)         ND (0.70)         ND (0.70)         ND (0.70)         ND (0.70)         ND (0.70)         ND (0.70)         ND (0.60)         ND (0.65	5			× /	× /	· · · · · ·	· /	~ /	``´´	· /
n-Propylbenzeneμg/L5ND (0.60)ND (0.60)ND (0.60)ND (0.60)ND (0.60)ND (0.60)1.9 J1.3 JStyreneμg/L5ND (0.70)ND (0.70)ND (0.70)ND (0.70)ND (0.70)ND (0.70)ND (0.70)ND (0.70)1,1,2Tetrachloroethaneμg/L5ND (0.60)ND (0.60)ND (0.60)ND (0.60)ND (0.60)ND (0.60)ND (0.60)ND (0.60)1,2,2-Tetrachloroethaneμg/L5ND (0.65)'ND (0.65)'ND (0.65)'ND (0.65)'ND (0.65)'ND (0.65)'ND (0.65)'ND (0.65)'Tetrachloroethaneμg/L5ND (0.65)'ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)Tolueneμg/L5ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.53)1,2,3-Trichlorobenzeneμg/L5ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)1,1,1-Trichloroethaneμg/L5ND (0.53)ND (0.54)ND (0.54)ND (0.54)ND (0.53)ND (0.53)1,1,2-Trichloroethaneμg/L5ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)1,1,2-Trichloroethaneμg/L5ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)1,1,2-Trichloroethaneμg/L5ND (0.53)I.2ND (0.53)ND (0.5				× /		· · · ·			× /	· /
Styrene $\mu g/L$ 5ND (0.70)ND (0.7	· ·				``´´	· · · · · ·	· /	``´´	, ,	· · · ·
1,1,2-Tetrachloroethane         µg/L         5         ND (0.60)         ND (0.65) <sup>c</sup> ND (0.50)         <				<b>`</b>	``´´	· · · ·	· · · ·	· · · · ·		
1,1,2,2-Tetrachloroethane $\mu g/L$ 5ND (0.65)°ND (0.50)ND (	-			<b>`</b>	, , , , , , , , , , , , , , , , , , ,	. ,	· · · ·		, , , , , , , , , , , , , , , , , , ,	· /
Tetrachloroetheneμg/L5ND (0.90)ND (0.90)ND (0.90)ND (0.90)ND (0.90)ND (0.90)ND (0.90)ND (0.90)Tolueneμg/L5ND (0.53)ND (0.53)ND (0.53)ND (0.53)0.66 JND (0.53)ND (0.53)1,2,3-Trichlorobenzeneμg/L5ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)1,2,4-Trichlorobenzeneμg/L5ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)1,1,1-Trichloroethaneμg/L5ND (0.54)ND (0.54)ND (0.54)ND (0.54)ND (0.54)ND (0.53)1,1,2-Trichloroethaneμg/L1ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)Trichloroethaneμg/L5ND (0.53)1.2ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)Trichloroethaneμg/L5ND (0.53)1.2ND (0.53)ND (0.53)ND (0.53)ND (0.53)Trichloroethaneμg/L5ND (0.53)1.2ND (0.53)ND (0.53)ND (0.53)ND (0.53)Trichloroethaneμg/L5ND (0.84)ND (0.84)ND (0.84)ND (0.84)ND (0.84)ND (0.84)1,2,3-Trichloropropaneμg/L0.04ND (0.70)ND (0.70)ND (0.70)ND (0.70)ND (0.70)				× /			. ,	× /	``´´	. ,
Tolueneμg/L5ND (0.53)ND (0.50)ND (0.53)ND (0.53)<				× /	× /	· /	``´´	× /	· /	. ,
1,2,3-Trichlorobenzene $\mu g/L$ 5ND (0.50)ND (0.50)<				· /	` <i>´</i>	· · · · · ·	· · · /		· · · · · · · · · · · · · · · · · · ·	
1,2,4-Trichlorobenzene $\mu g/L$ 5ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)ND (0.50)1,1,1-Trichloroethane $\mu g/L$ 5ND (0.54)ND (0.54)ND (0.54)ND (0.54)ND (0.54)ND (0.54)ND (0.54)1,1,2-Trichloroethane $\mu g/L$ 1ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)Trichloroethene $\mu g/L$ 5ND (0.53)1.2ND (0.53)ND (0.53)ND (0.53)ND (0.53)Trichlorofluoromethane $\mu g/L$ 5ND (0.84)ND (0.84)ND (0.84)ND (0.84)ND (0.84)ND (0.84)1,2,3-Trichloropropane $\mu g/L$ 0.04ND (0.70)ND (0.70)ND (0.70)ND (0.70)ND (0.70)				<b>`</b>	``´´	· · · ·	. ,		( )	· · · · ·
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1,1,2-Trichloroethane $\mu g/L$ 1ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)Trichloroethene $\mu g/L$ 5ND (0.53)1.2ND (0.53)ND (0.53)ND (0.53)ND (0.53)ND (0.53)Trichlorofluoromethane $\mu g/L$ 5ND (0.84)ND (0.84)ND (0.84)ND (0.84)ND (0.84)ND (0.84)ND (0.84)1,2,3-Trichloropropane $\mu g/L$ 0.04ND (0.70)ND (0.70)ND (0.70)ND (0.70)ND (0.70)				<b>`</b>			, ,		. ,	
Trichloroethene $\mu g/L$ 5         ND (0.53)         1.2         ND (0.53)         ND (0.54)         ND (0.54)         ND (0.54)         ND (0.70)         ND (0.7	1,1,1-Trichloroethane		5							
Trichlorofluoromethane         μg/L         5         ND (0.84)         ND (0.70)         <			1	<b>`</b>	, <i>,</i> ,	. ,			```´	
1,2,3-Trichloropropane µg/L 0.04 ND (0.70)	Trichloroethene			· /					. ,	
				. ,					· · · ·	
1,2,4-Trimethylbenzene $\mu g/L$ 5         ND (1.0)         ND (1.0)         ND (1.0)         ND (1.0)         ND (1.0)         ND (1.0)		µg/L			. ,	. ,	· /	· · · ·	<b>`</b>	. ,
	1,2,4-Trimethylbenzene	μg/L	5	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)

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#### TABLE 2 - SUMMARY OF GROUNDWATER ANALYTICAL RESULTS - VOCs and PCBs

Date Sampled: December 5, 2018

Client Sample ID:	Unit	State	MW-1-D	<b>MW-2</b>	MW-5	MW-6	MW-9	MW-10	MW-12
Lab Sample ID:	Unit	<b>Standard</b> <sup>1</sup>	JC79316-6	JC79316-5	JC79316-4	JC79316-3	JC79316-1	JC79316-8	JC79316-7
			MS Vola	tiles (SW84	6 8260C)				
1,3,5-Trimethylbenzene	μg/L	5	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)	ND (1.0)
Vinyl chloride	μg/L		24	ND (0.79)	ND (0.79)	ND (0.79)	ND (0.79)	ND (0.79)	ND (0.79)
m,p-Xylene	μg/L	-	ND (0.78)	ND (0.78)	ND (0.78)	ND (0.78)	ND (0.78)	ND (0.78)	ND (0.78)
o-Xylene	μg/L	5	ND (0.59)	ND (0.59)	ND (0.59)	ND (0.59)	ND (0.59)	ND (0.59)	ND (0.59)
TOTAL VOCs	μg/L		29.48	17.69	18.22	35.81	11.06	13.06	24.42
			GC/LC Semi	-volatiles (S	W846 8082A	.)			
Aroclor 1016	μg/L	0.09	ND (0.094)	ND (0.094)	ND (0.094)	ND (0.094)	ND (0.094)	ND (0.094)	ND (0.098)
Aroclor 1221	μg/L	0.09	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.20)	ND (0.21)
Aroclor 1232	μg/L	0.09	ND (0.12)	ND (0.12)	ND (0.12)	ND (0.12)	ND (0.12)	ND (0.12)	ND (0.13)
Aroclor 1242	μg/L	0.09	ND (0.11)	ND (0.11)	ND (0.11)	ND (0.11)	ND (0.11)	ND (0.11)	ND (0.11)
Aroclor 1248	μg/L	0.09	ND (0.061)	<b>0.75</b> <sup>e</sup>	ND (0.061)	ND (0.061)	ND (0.061)	ND (0.061)	ND (0.063)
Aroclor 1254	μg/L		ND (0.20)	1.1	ND (0.20)	ND (0.20)	ND (0.20)	0.28	ND (0.21)
Aroclor 1260	μg/L		ND (0.073)	1.2	ND (0.073)	ND (0.073)	ND (0.073)	ND (0.073)	ND (0.076)
TOTAL PCBs	μg/L	0.09	ND	3.05	ND	ND	ND	0.28	ND

Notes:

Legend: Hit Exceed

<sup>1</sup>DEC Division of Water Technical and Operational Guidance Series (TOGS) 1.1.1,

Ambient Water Quality Standards and Guidance Values, reissued June 1998.

\*State standard is 5  $\mu$ g/L for each xylene isomer.

<sup>a</sup>Associated CCV outside of control limits low.

<sup>b</sup>Associated CCV outside of control limits high, sample was ND.

°This compound in BS is outside in house QC limits bias high.

<sup>d</sup>Associated CCV outside of control limits high, sample was ND. This compound in BS is outside in house QC limits bias high.

<sup>e</sup>More than 40 % RPD for detected concentrations between the two GC columns.

 $\mu$ g/L micrograms per liter, equivalent to parts per billion (ppb)

ND Not Detected

--- No promulgated State Standard

J Indicates an estimated value

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	State										Mor	nitoring V	Well Loca	ation									
Compound	Standard <sup>1</sup>	MW	V-1S			MW	/-1D					<b>MW-2</b>			M	V-3	M	W-4			MW-5		
	(µg/L)	03/12/08	09/23/15	03/12/08	01/05/09	09/23/15	12/16/16	10/05/17	12/05/18	03/12/08	09/23/15	12/16/16	10/05/17	12/05/18	03/12/08	09/23/15	03/12/08	09/23/15	03/12/08	09/23/15	12/16/16	10/05/17	12/05/18
Acetone	50	NS	NS		20	13,200				NS		7											
1,1,2-Trichloroethane	1	NS	NS							NS													
1,1-Dichloroethane	5	NS	NS		2		2	2.9	2.3	NS		3		0.73 J									
1,2,3-Trimethylbenzene	5	NS	NS		2				NA	NS				NA									NA
1,2,4-Trichlorobenzene	5	NS	NS							NS		1											
1,2,4-Trimethylbenzene	5	NS	NS		6.0					NS													
1,2-Dibromo-3-Chloropropane	0.04	NS	NS							NS													
1,2-Dichlorobenzene	3	NS	NS		2.0		0.5	0.55 J	0.59 J	NS		1.5	3.2	3.1									
1,2-Dichloroethene (Total)	5	NS	NS		2.0		1.6		1.5	NS		1.4	2	1									
1,2-Dichloropropane	1	NS	NS							NS													
1,3,5-Trimethylbenzene	5	NS	NS		2.0					NS													
1,3-Dichlorobenzene	3	NS	NS							NS		2.8	3.1	2.5									
1,4-Dichlorobenzene	3	NS	NS							NS		5.9	6.6	5.9									
2-Butanone (MEK)	50	NS	NS	52,000	7,800	10,100				NS													
Benzene	1	NS	NS		2.0		0.5	0.83	0.52	NS	0.8	0.9	3	1.4									
Bromodichloromethane	50	NS	NS	NA	NA	NA	NA	NA		NS	NA	NA	NA		NA	0.82 J							
Carbon Disulfide	60	NS	NS				0.4			NS			0.50 J										
Chlorobenzene	5	NS	NS		2.0		0.5	0.49 J	0.57 J	NS	1.9	1.2	5.5	0.97 J									
Chloroethane	5	NS	NS				0.7			NS		1.3	2.1	0.89 J									
Chloroform	7	NS	NS							NS											3.2	15.4	17.4
Ethylbenzene	5	NS	NS		3.0					NS		0.2	0.27 J										
Isopropylbenzene	5	NS	NS		1.0					NS			0.60 J										
m/p-Xylenes	5	NS	NS		5.0					NS													
Methyl tert-butyl ether	10	NS	NS		0.5		0.4	0.55 J		NS													
Methylene Chloride	5	NS	NS							NS													
n-Butylbenzene	5	NS	NS							NS		0.9											
n-propylbenzene	5	NS	NS		2.0					NS		0.2											
o-Xylene	5	NS	NS							NS													
p-Isopropyltoluene	5	NS	NS							NS													
sec-Butylbenzene	5	NS	NS		-					NS													
tert-Butylbenzene	5	NS	NS							NS													
Toluene	5	NS	NS		0.9					NS		0.3	0.30 J										
Tetrachloroethene	5	NS	NS							NS	1.8												
Trichloroethene	5	NS	NS							NS	2.5	5.0	1.5	1.2									
Vinyl Chloride	2	NS	NS		23.0	2.3	28.5	13.4	24	NS		0.9											
Total VOCs		NS	NS	52,000	7,875	23,302	36	17	28	0	7	33	27	15	0	0	0	0	0	0	3	15	18

### TABLE 3 - SUMMARY OF HISTORICAL GROUNDWATER ANALYTICAL RESULTS - VOCs [DETECTIONS ONLY] (µg/L)

Plumley Engineering, P.C.

Project No. 2015127

# TABLE 3 - SUMMARY OF HISTORICAL GROUNDWATER ANALYTICAL RESULTS - VOCs [DETECTIONS ONLY

	State										Monitor	ing Well	Location									
Compound	Standard <sup>1</sup>			MW-6			MV	<i>N</i> -7	MV	W-8			MW-9					MW-10			MV	V-11
	(µg/L)	01/05/09	09/23/15	12/16/16	10/05/17	12/05/18	01/05/09	09/23/15	01/05/09	09/23/15	01/05/09	09/23/15	12/16/16	10/05/17	12/05/18	01/05/09	09/23/15	12/16/16	10/05/17	12/05/18	06/29/09	09/23/15
Acetone	50			22			NS											8				
1,1,2-Trichloroethane	1																	2				
1,1-Dichloroethane	5	2	1	1	1.4	0.61 J	NS									4		0.4	0.64 J		2	
1,2,3-Trimethylbenzene	5					NA	NS				2				NA	22				NA		
1,2,4-Trichlorobenzene	5						NS									1		1				
1,2,4-Trimethylbenzene	5						NS							3.1		100.0		4.3	3.1		2.0	
1,2-Dibromo-3-Chloropropane	0.04						NS									2.0						
1,2-Dichlorobenzene	3	9.0	4.8	3.7	7.1	4.5	NS	1.8	2.0		5.0	6.7	1.9	3.3	1	18.0	10.7	6.3	11.4	2.6	3.0	2.8
1,2-Dichloroethene (Total)	5	1.0	1.1	0.5			NS											1.2	0.77	0.93 J	4.0	
1,2-Dichloropropane	1						NS				1.0											
1,3,5-Trimethylbenzene	5						NS						0.4	1.0 J		25.0		0.7	0.50 J			
1,3-Dichlorobenzene	3	5.0	2.7	1.6	3.2	2.4	NS				2.0	1.9	0.9	0.99 J	0.62 J			0.5	0.68 J			
1,4-Dichlorobenzene	3	4.0	2.7	1.8	3.8	2.7	NS	1.2			4.0	4.4	2.2	2.3	1.6	3.0	2.9	1.7	3.5	1.1		
2-Butanone (MEK)	50						NS															
Benzene	1	3.0	2.1	1.8	3	1.7	NS				4.0	2.2	0.4	1.2	0.66	41.0	25.6	3.1	24.2	2.3	0.6	
Bromodichloromethane	50	NA	NA	NA	NA		NS	NA	NA	NA	NA	NA	NA	NA		NA	NA	NA	NA		NA	NA
Carbon Disulfide	60						NS															
Chlorobenzene	5	48.0	21.0	16.4	33.6	23.9	NS				30.0	29.1	7.8	14.1	5.8	3.0	2.0	0.5	2.2		4.0	3.1
Chloroethane	5						NS									12.0	4.1		2.3			
Chloroform	7						NS		0.8		2.0											
Ethylbenzene	5						NS							1.8		51.0		0.4	0.65 J			
Isopropylbenzene	5	5.0		0.3	0.27 J		NS		3.0		9.0	11.1	1.9	6.6	0.72 J	23.0	9.7	6.1	12.6	2.3		
m/p-Xylenes	5						NS									83.0	2.0	1.8			1.0	
Methyl tert-butyl ether	10						NS															
Methylene Chloride	5						NS				5.0											
n-Butylbenzene	5	1.0					NS		2.0		5.0		0.4	3.3		5.0		2.4	2	0.63 J		
n-propylbenzene	5	4.0					NS		2.0		12.0	11.7	1.2	7.1		26.0	10.3	9.0	12.2	1.9 J		
o-Xylene	5						NS				1.0			0.52 J		11.0	1.2	0.3	1.3		0.9	
p-Isopropyltoluene	5						NS							0.26		4.0						
sec-Butylbenzene	5	2.0			0.39 J		NS		3.0		4.0		1.1	3.3		5.0		2.5	3.3	1.3 J		
tert-Butylbenzene	5	1.0			0.45 J		NS						0.4	0.81 J		1.0		0.7	0.90 J			
Toluene	5						NS				3.0	2.2	0.4	1	0.66 J	16.0	2.1	0.7	2.6			
Tetrachloroethene	5			0.2			NS															
Trichloroethene	5			0.3			NS	1.1								2.0		1.6	0.50 J			1.4
Vinyl Chloride	2	5.0	2.2	0.8	0.62 J		NS															
Total VOCs		90	38	50	52	36	NS	4	13	0	89	69	19	47	11	458	71	54	81	13	18	7

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<b>Y</b> ]	(µg/L)
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Project No. 2015127

#### **Monitoring Well Location** State **MW-12** Compound Standard<sup>1</sup> **RW-1 RW-2** RW-3 **RW-4** Notes: 07/14/09 $(\mu g/L)$ 12/05/18 09/23/15 09/23/15 12/16/16 10/05/17 09/23/15 09/23/15 09/23/15 7.0 50 Acetone --------------------------1,1,2-Trichloroethane 1 --------------------------5 VOCs 1,1-Dichloroethane 2.3 ------------------------1,2,3-Trimethylbenzene 5 PCBs ---NA ----------------------1,2,4-Trichlorobenzene 5 NA ----------------------------1,2,4-Trimethylbenzene 5 NS ---------------------------1,2-Dibromo-3-Chloropropane 0.04 μg/L ----------------------------1,2-Dichlorobenzene 3 6.0 6.4 4.8 3.3 3.4 3.4 4.6 5.6 5.6 ---1,2-Dichloroethene (Total) 5 -------------------------------1,2-Dichloropropane 1 ---------------------------1,3,5-Trimethylbenzene 5 ----------------------------1,3-Dichlorobenzene 3 0.3 3.7 ---------------------1,4-Dichlorobenzene 3 1.3 1.7 1.2 1.8 7.1 3.2 1.8 2.0 ---2-Butanone (MEK) 50 ------------------------------1.0 0.5 0.20 J 0.46 J 3.7 2.5 1.3 Benzene 1 ---1.1 Bromodichloromethane 50 NA NA NA NA NA NA NA NA ---Carbon Disulfide 60 -----------------------------Chlorobenzene 5 9.1 4.3 1.8 1.8 1.1 ---10.1 12.1 11.2 5 2.5 Chloroethane 2.0 ----------------------Chloroform 7 ------------------------------5 Ethylbenzene ----------------------------Isopropylbenzene 5 1.0 ---1.8 ---3.9 -----------m/p-Xylenes 5 ---------------------------Methyl tert-butyl ether 10 ---------------------------Methylene Chloride 5 --------------------------n-Butylbenzene 5 --------------------------n-propylbenzene 5 1.3 J 0.4 --------------------o-Xylene 5 ----------------------------p-Isopropyltoluene 5 --------------------------sec-Butylbenzene 5 ---0.51 J 1.4 J -----------------tert-Butylbenzene 5 0.88 J 0.96 J 0.9 ------------------Toluene 5 -----------------------------5 Tetrachloroethene ----------------------------Trichloroethene 5 1.7 1.7 ----------------------Vinyl Chloride 2 ---------------------------Total VOCs 17 17 21 17 24 26 15 12 ---10

#### TABLE 3 - SUMMARY OF HISTORICAL GROUNDWATER ANALYTICAL RESULTS - VOCs [DETECTIONS ONLY] (µg/L)

Volatile Organic Compounds Polychlorinated Biphenyls Not Analyzed

Not Sampled

micrograms per liter, equivalent to parts per billion (ppb) Indicates the specified compound was not detected at a concentration exceeding the method detection limit.

Plumley Engineering, P.C.

<sup>1</sup>DEC Division of Water's Technical and Operational Guidance Series (TOGS) 1.1.1, Ambient Water Quality Standards and Guidance Values, reissued June 1998.

#### TABLE 4 - SUMMARY OF HISTORICAL GROUNDWATER ANALYTICAL RESULTS - TOTAL PCBs (µg/L)

Commoniad	State Standard <sup>1</sup>	Sample								Monitor	ring Well l	Location							
Compound	(µg/L)	Date	MW-1D	MW-1S	<b>MW-2</b>	<b>MW-3</b>	MW-4	<b>MW-5</b>	<b>MW-6</b>	<b>MW-7</b>	<b>MW-8</b>	MW-9	<b>MW-10</b>	<b>MW-11</b>	<b>MW-12</b>	<b>RW-1</b>	<b>RW-2</b>	<b>RW-3</b>	RW-4
Total PCBs	0.09	03/12/2008		FP	FP				NS	NI	NI	NI	NI	NI	NI	NI	NI	NI	NI
Total PCBs	0.09	2009*		FP	FP	NS	NS	NS		FP	0.13	0.93				NI	NI	NI	NI
		12/08/2011							R	emedial E	Excavation	Complete	ed						
Total PCBs	0.09	09/23/2015	0.32	FP	2.5					11.3			0.29			5.5	3.3	0.23	
Total PCBs	0.09	12/16/2016		FP	12	NS	NS			FP	NS		2.9	NS		NS	NS	NS	NS
Total PCBs	0.09	10/05/2017		FP	4.3	NS	NS			FP	NS			NS		NS	NS	NS	NS
Total PCBs	0.09	12/05/2018		NS	3.05	NS	NS			NS	NS		0.28	NS		NS	NS	NS	NS
																T	T*4		
Notes:															Legend:	H	lit	Exc	ee:

<sup>1</sup>DEC Division of Water's Technical and Operational Guidance Series (TOGS) 1.1.1, Ambient Water Quality Standards and Guidance Values, reissued June 1998. State standard for PCBs is 0.09  $\mu$ g/L for each Aroclor.

- PCBs Polychlorinated Bipenyls
- μg/L Micrograms per liter
- NI Not Installed
- NS Not Sampled
- Free Product Present Not Sampled FP

Indicates the specified compound was not detected at a concentration exceeding the method detection limit. ---

\*2009 samples collected on 1/5/2009, 6/29/2009 and 7/14/2009

Refer to laboratory reports for additional information.

#### TABLE 5 - PFAS AND 1,4-DIOXANE IN GROUNDWATER

Date Sampled: December 5, 2018

Client Sample ID:	Unit	State	<b>MW-9</b>	EQUIPMENT BLANK	MW-12	MW-10
Lab Sample ID:	- •	Standard <sup>1</sup>	JC79316-1	JC79316-2	JC79316-7	JC79316-8
I	MS Sei	mi-volatiles	(EPA 537M	BY ID)		
Perfluorobutanoic acid	ng/L	-	10.5 J	ND (1.9)	4.87 J	ND (38)
Perfluoropentanoic acid	ng/L	-	ND (2.9)	ND (1.4)	ND (1.5)	ND (29)
Perfluorohexanoic acid	ng/L	-	7.03 J	ND (0.96)	1.31 J	ND (19)
Perfluoroheptanoic acid	ng/L	-	2.09 J	ND (0.96)	ND (1.0)	ND (19)
Perfluorooctanoic acid	ng/L	-	23.6ª	ND (0.96)	3.41	ND (19) <sup>b</sup>
Perfluorononanoic acid	ng/L	70	ND (1.9)	ND (0.96)	ND (1.0)	ND (19)
Perfluorodecanoic acid	ng/L	-	ND (1.9)	ND (0.96)	ND (1.0)	ND (19)
Perfluoroundecanoic acid	ng/L	-	ND (1.9)	ND (0.96)	ND (1.0)	ND (19)
Perfluorododecanoic acid	ng/L	-	ND (2.9)	ND (1.4)	ND (1.5)	ND (29)
Perfluorotridecanoic acid	ng/L	-	ND (1.9)	ND (0.96)	ND (1.0)	ND (19)
Perfluorotetradecanoic acid	ng/L	-	ND (1.9)	ND (0.96)	ND (1.0)	ND (19)
Perfluorobutanesulfonic acid	ng/L	-	1.93 J	ND (0.96)	ND (1.0)	ND (19)
Perfluorohexanesulfonic acid	ng/L	-	2.87 J	ND (0.96)	ND (1.0)	ND (19)
Perfluoroheptanesulfonic acid	ng/L	-	ND (1.9)	ND (0.96)	ND (1.0)	ND (19)
Perfluorooctanesulfonic acid	ng/L	-	3.40 J	ND (1.4)	6.12	ND (29)
Perfluorodecanesulfonic acid	ng/L	-	ND (1.9)	ND (0.96)	ND (1.0)	ND (19)
PFOSA	ng/L	-	ND (1.9)	ND (0.96)	ND (1.0)	ND (19)
MeFOSAA	ng/L	-	ND (7.7)	ND (3.8)	ND (4.0)	ND (77)
EtFOSAA	ng/L	-	ND (7.7)	ND (3.8)	ND (4.0)	ND (77)
6:2 Fluorotelomer sulfonate	ng/L	-	ND (3.8)	ND (1.9)	ND (2.0)	ND (38)
8:2 Fluorotelomer sulfonate	ng/L	-	ND (3.8)	ND (1.9)	ND (2.0)	ND (38)
MS	5 Semi	-volatiles (S	W846 8270D	BY SIM)		
1,4-Dioxane	μg/L	-	2.34	-	0.239	0.225

Notes:

Legend: Hit Exceed

<sup>1</sup>DEC Division of Water Technical and Operational Guidance Series (TOGS) 1.1.1, *Ambient Water Quality Standards and Guidance Values*, reissued June 1998.

<sup>a</sup>Associated CCV outside of control limits high.

<sup>b</sup>Associated CCV outside of control limits high, sample was ND.

PFAS Per- and Poly-Fluorinated Alkyl Substances

- ng/L nanograms per liter, equivalent to parts per trillion (ppt)
- $\mu$ g/L micrograms per liter, equivalent to parts per billion (ppb)

ND Not Detected

J Indicates an estimated value

#### 2802-2810 LODI STREET

# DEC Site No. 7-34-013

City of Syracuse, Onondaga County, New York

Inspector:

Company:

un

Recovery Well	Well Head Conditions: OK / Not OK	Depth to Water (feet)	Free Product: Present (Yes) Absent (No)	Free Product Thickness (inches)	Free Product Volume Removed (gallons)	Comments
MW-1S	OK	2.680	Y	1.75	2 Full	Replan
MW-2	( -	30.41	~			
MW-7		23.95	Y	0.3	1/2	Roolun
MW-10		26.90	$\sim$			
RW-1		28.06	N			
RW-2		26.61	N		0	
RW-3		27,98	YN	-	1/4	
RW-4	V	27.15	N			

Monitoring Well	Well Head Conditions: OK / Not OK	Depth to Water (feet)	Free Product: Present (Yes) Absent (No)	Comments
MW-1D			11000110 (110)	
MW-3				
MW-4				
MW-5				
MW-6				
MW-8				
MW-9				
MW-11				
MW-12				

Plumley Engineering, P.C.

2/28/18

Date:

#### 2802-2810 LODI STREET

**DEC Site No. 7-34-013** 

City of Syracuse, Onondaga County, New York

Inspector:	MTM	
Company:	glu	in h

5/30/18 Date:

		<u>,,,,,</u>						
Recovery Well	Well Head Conditions: OK / Not OK	Depth to Water (feet)	Free Product: Present (Yes) Absent (No)	Free Product Thickness (inches)	Free Product Volume Removed (gallons)	Comments		
MW-1S	on	29.11	Ý	0.8'fr	٢ '/4	Replace Sochs		
MW-2		21.55	N			V J		
MW-7		25.62	Y	6.1	f.Im			
MW-10		27.99	4	6.1	f.Ir	×		
RW-1		26.89	N			Nort		
RW-2		27.33	1					
RW-3		26.72						
RW-4	17	25.87	P P			U		
				тт				
Monitoring Well	Well Head Conditions: OK / Not OK	Depth to Water (feet)	Free Product: Present (Yes) Absent (No)		Comm	ents		
-	Well Head Conditions:	Water (feet)	Present (Yes)		Comm	ents		
Well	Well Head Conditions:	Water (feet) 31,95	Present (Yes) Absent (No)		Comm	ents		
Well MW-1D	Well Head Conditions: OK / Not OK	Water (feet)	Present (Yes) Absent (No)		Comm	ents		
Well MW-1D MW-3	Well Head Conditions: OK / Not OK	Water (feet) 31,95	Present (Yes) Absent (No)		Comm	ents		
Well MW-1D MW-3 MW-4	Well Head Conditions: OK / Not OK	Water (feet) 31,95 25.91 	Present (Yes) Absent (No)		Comm	ents		
Well MW-1D MW-3 MW-4 MW-5	Well Head Conditions: OK / Not OK	Water (feet) 31,95 25,91 	Present (Yes) Absent (No)		Comm	ents		
Well MW-1D MW-3 MW-4 MW-5 MW-6	Well Head Conditions: OK / Not OK	Water (feet) 31,95 25.91 	Present (Yes) Absent (No)		Comm	ents		
Well           MW-1D           MW-3           MW-4           MW-5           MW-6           MW-8	Well Head Conditions: OK / Not OK	Water (feet) 31,95 25.91 	Present (Yes) Absent (No)		Comm	ents		

Plumley Engineering, P.C.

Project No. 2010131

#### 2802-2810 LODI STREET

**DEC Site No. 7-34-013** 

City of Syracuse, Onondaga County, New York

Inspector:
------------

Company:

TN

Plumber

Recovery Well	Well Head Conditions: OK / Not OK	Depth to Water (feet)	Free Product: Present (Yes) Absent (No)	Free Product Thickness (inches)	Free Product Volume Removed (gallons)	Comments
MW-1S		28.69	Y	0.3'54	< 1/4	Repare Such
MW-2		26.89	N	N		0
MW-7		24.97				
MW-10		27.19				
RW-1		26.08				
RW-2		26.07				
RW-3		26.01				
RW-4		25.17	1	P		

Monitoring	Well Head	Depth to	Free Product:	
Well	<b>Conditions:</b>	Water	Present (Yes)	Comments
wen	OK / Not OK	(feet)	Absent (No)	
MW-1D				
MW-3				
MW-4				
MW-5				
MW-6				
MW-8				
MW-9				
MW-11			· · · · · · · · · · · · · · · · · · ·	
MW-12				

Plumley Engineering, P.C.

9/28/18

Date:

#### 2802-2810 LODI STREET

**DEC Site No. 7-34-013** 

City of Syracuse, Onondaga County, New York

Inspector:	NTM	
Company:	Plumby	

Recovery Well	Well Head Conditions: OK / Not OK	Depth to Water (feet)	Free Product: Present (Yes) Absent (No)	Free Product Thickness (inches)	Free Product Volume Removed (gallons)	Comments
MW-1S	OK	29.75	Ý	4.158	21/2 Cullu	Replaced /
MW-2	1	27.72				Lemsrod
MW-7		23.86	Ý	0.1	1 sock	Replach
MW-10		24,09	Y	6.10		Remarch
RW-1		27.32	·			
RW-2		26.95				
RW-3		26.40	Ý	<.40		Remarch
RW-4	$\checkmark$	25.41				

Monitoring Well	Well Head Conditions: OK / Not OK	Depth to Water (feet)	Free Product: Present (Yes) Absent (No)	Comments
MW-1D	OK	30.55	N	
MW-3	(	25.40	1	
MW-4	Cao	26.24		
MW-5	Can	22.25		
MW-6	Cao	29.97	l l	
MW-8	1	26.41		
MW-9		17,59		
MW-11	4	29,17	1 1	
MW-12	(40	26.80	↓/	

Plumley Engineering, P.C.

12/5/18

Date:

# GROUNDWATER SAMPLING FIELD LOGS

Client/Site:	Quanta Resources	l		Project No.:	2015007
<b>Monitoring Location:</b>				_ Date:	12/5/2018
Source Description:	MW-1D			_ Sampler:	MTM/DTH
Well & Water Level D	Data:	Tota	l Depth of Well	50.19	feet
		<b>Initial</b>	Depth to Water	30,55	feet
	I	Length of Water	Column (LWC)		feet
Purge Volume Calcula	ation:				
Well Diameter (in	nches):	<b>Calculated W</b>	ell Volume To l	Be Removed	
1	,	LWC * 0.041		Gallons	
1.25		LWC * 0.064	* 3 =	Gallons	
1.5		LWC * 0.092		Gallons	
2		LWC * 0.163	* 3 = 10	Gallons	
3		LWC * 0.367		Gallons	
4		LWC * 0.653		Gallons	
6		LWC * 1.469		Gallons	
Free Product Check:	Free P	Product Present:	Yes	No	
	Measured Thick		105		
Purge Data:	Purge Date:	-			
				_	
	<b>Purging Time:</b>	From:	12:15	- То:	12.30
1	Type of Purging E	quipment Used:	Whale F	ump	
	Purged Wa	ater Comments:		br - noslea	4
Sampling Data:	Depth to Wat	ter at Sampling:	30.2		feet
	•				ieet
	Color of Sample:	cleve	Sample Date:		
	Turbidity:_	-	Sample Time	14.00	
Ту	pe of Sampling E	quipment Used:	Bu	hr	
<b>Field Indicators P</b>	resent During San	nple Collection:	Odor		
	0	-	Sheen		
			<b>Free Product</b>	1	
			None	×	
Notes:				2	
X					
			$\cap$		
Weather:	Temperature <sup>o</sup> F	40	Sunny Cloudy	Rain Snow	
		1			

Client/Site:	Quanta Resources				<b>Project No.:</b>	2015007
<b>Monitoring Location:</b>	-				Date:	12/5/2018
Source Description:	MW-1S				Sampler:	MTM/DTH
Well & Water Level D	Data:	Total	Depth	of Well:	38.9	feet
					500	feet
	Le	ength of Water C				feet
Purge Volume Calcula	ation.					
Well Diameter (in		Calculated We	all Volu	me To Re	Removed	
1	nenes).	LWC * 0.041				
1.25		LWC * 0.064			- Gallons	
1.23		LWC * 0.092			- Gallons	
2		LWC * 0.163			-	
3		LWC * 0.367				
4		LWC * 0.653		-	- Gallons	
6		LWC * 1.469			- Gallons	
			1			
Free Product Check:		oduct Present:	Y	es	No	
	Measured Thickn	ess/Comment:		4,7	15	
Purge Data:	Purge Date:					
	<b>Purging Time:</b>				То:	
1	Type of Purging Eq	uipment Used:				
	Purged Wat	ter Comments:				
Sampling Data:	Depth to Wate	r at Sampling: _				feet
(	Color of Sample:		Samn	le Date:		
	Turbidity:		-	e Time:		
т.	•	uinmont Used.		-		ai
1 y	pe of Sampling Eq	ulpment Used:				
<b>Field Indicators P</b>	resent During Sam	ple Collection:	Od	or		
			She	en		
			Free P	roduct		
			No	ne		
Notes:				1		
19		(No	SU	mple	)	
·	2 0.10 year 1	remond				
	0					
Weather:	Temperature <sup>o</sup> F		lunnv	Cloudy	Rain Snow	
		<b>`</b>	J	Juna		

Client/Site:	Quanta Resources			Project No.:	
<b>Monitoring Location:</b>	·			Date:	
Source Description:	<u>MW-2</u>			Sampler:	MTM/DTH
Well & Water Level I	Data:	Tota	al Depth of We	ell: 37.9	feet
		Initial	Depth to Wate	er: 27.70	feet
	Le	ength of Water	Column (LWC	C):	feet
Purge Volume Calcula	ation:				
Well Diameter (i		Calculated W	Vell Volume To	Be Removed	
1	,	LWC * 0.041		Gallons	
1.25		LWC * 0.064	<b>1 * 3</b> =	Gallons	
1.5		LWC * 0.092	2 * 3 =	Gallons	
2		LWC * 0.163	3 * 3 = 5	Gallons	
3		LWC * 0.367	7 * 3 =	Gallons	
4		LWC * 0.653		Gallons	
6		LWC * 1.469	77	Gallons	
Free Product Check:	Free Pr	oduct Present:	Yes	No	
	Measured Thickn			2	
Purge Data:	Purge Date:	12	5/18		
	<b>Purging Time:</b>	From:	12:00	То:	12:10
ſ	<b>Fype of Purging Eq</b>	uipment Used:	Whi	le Pura	
	Purged Wat	ter Comments:	sligh		<
Sampling Data:	Depth to Wate	r at Sampling:	27.91		feet
	Color of Sample:	chur	Sample Dat	e: 12/5/18	
	Turbidity:	sligint	Sample Tim		e 
Ту	pe of Sampling Eq		B	sa. li (	
<b>Field Indicators P</b>	resent During Sam	ple Collection:	Odor	×	
		-	Sheen		2
			Free Produc	t	5
			None		3
Notes:					
			0		
Weather:	Temperature <sup>o</sup> F	10	Sunny Cloud	ly Rain Snow	

Client/Site: Monitoring Location:	Quanta Resources			Project No.: Date:	2015007
Source Description:	MW-5			Sampler: MT	
Well & Water Level I Purge Volume Calcula	Data: Length ation:	Initial De of Water Co	Depth of Well: epth to Water: olumn (LWC):	42.01 fee 27.25 fee	t t
Well Diameter (i	· · · · · · · · · · · · · · · · · · ·	lculated Wel VC * 0.041 *	<u>ll Volume To Be</u> 2 –		
1 1.25		VC * 0.041 * VC * 0.064 *		_ Gallons Gallons	
1.23		VC * 0.092 *		_ Gallons	
2		VC * 0.163 *		_ Gallons	
3		VC * 0.367 *	0	Gallons	
4	LW	VC * 0.653 *	3 =	Gallons	
6	LW	VC * 1.469 *	3 =	Gallons	
Free Product Check:	Free Product Measured Thickness/C		Yes	No	
Purge Data:	Purge Date:	12/5			
	<b>Purging Time:</b>	From:	11:45	То:	1132
ן	Type of Purging Equipm Purged Water Co	S	B Whale oker	Purp	
Sampling Data:	Depth to Water at S	ampling:	27.3(	) feet	:
	Color of Sample: Turbidity:	LOT	Sample Date: _ Sample Time: _	12/14	
Ту	pe of Sampling Equipme	ent Used:	Bailes		
Field Indicators P	resent During Sample C		Odor Sheen Free Product		
Notes:			None	X	
Weather:	Temperature <sup>0</sup> F	10 6	unny Cloudy	Rain Snow	

Client/Site:	Quanta Resources			<b>Project No.:</b>	2015007
<b>Monitoring Location:</b>	oring Location:				12/5/2018
Source Description:	MW-6			Sampler:	MTM/DTH
Well & Water Level I	Data:		l Depth of Well:	41.49	-3
			Depth to Water:	29.91	feet
	Le	ength of Water (	Column (LWC):	11.5 4	feet
Purge Volume Calcul	ation:				
Well Diameter (i	nches):	Calculated W	ell Volume To B	e Removed	
1		LWC * 0.041	* 3 =	Gallons	
1.25		LWC * 0.064	* 3 =	Gallons	
1.5		LWC * 0.092	* 3 =	Gallons	
2		LWC * 0.163	* 3 = 10	Gallons	
3		LWC * 0.367	*3 =	Gallons	
4		LWC * 0.653	* 3 =	Gallons	
6		LWC * 1.469	* 3 =	Gallons	
Free Product Check:	Free Pr	oduct Present:	Yes	No	
The House check.	Measured Thickn		105	Ce	
D D t			1. 1		
Purge Data:	Purge Date:	120	10	ė	
	Purging Time:	From:	11:30	То:	11:40
r	<b>Fype of Purging Eq</b>	uipment Used:	What	Pump	
	Purged Wat	ter Comments:	Pink	shale	
Sampling Data:	Depth to Wate	r at Sampling:	30	03	feet
	Color of Sample:	pinkish	Sample Date:	115/11	
	Turbidity:	Stun	Sample Time:	13-01	•
T.		uinment Used.	Build		-3
1 )	vpe of Sampling Eq	urpment Useu:	DUU		
Field Indicators <b>P</b>	resent During Sam	ple Collection:	Odor		<b>4</b> 0
			Sheen	ie	-0
			Free Product	100.00	• 5
Notes:			None	X	el construction de la constructi
Notes:					
			~		
Weather:	Temperature <sup>0</sup> F	to	Sunny Cloudy	Rain Snow	
Revised 08/15/07					

Client/Site:	Quanta Resources			Project No.: Date:	2015007	
Monitoring Location Source Description:					MTM/DTH	
Well & Water Level	Data:	Initial	l Depth of Well: Depth to Water: Column (LWC):	36.53 feet		
Purge Volume Calcu Well Diameter ( 1 1.25 1.5 2 3		<u>Calculated W</u> LWC * 0.041 LWC * 0.064 LWC * 0.092 LWC * 0.163 LWC * 0.367	$\begin{array}{c} *3 = \\ *3 = \\ *3 = \\ 43 = 44 \end{array}$	<u>e Removed</u> Gallons Gallons Gallons Gallons Gallons		
4		LWC * 0.653	S.	Gallons		
6 Free Product Check: Purge Data:	Free P Measured Thick Purge Date:	8	Yes NA	Gallons		
Turge Data.	-		1.00 19 2	_		
а ж	Purging Time: Type of Purging Ec Purged Wa	From: uipment Used: ter Comments:	1070 Bluds, Pun cleor	-		
Sampling Data:	Depth to Wate	er at Sampling:			feet	
	Color of Sample: Turbidity: Type of Sampling Eq	slight	Sample Date: Sample Time: Blader	12.30	, Aur	
Field Indicators	Present During San	ple Collection:	Odor Sheen Free Product None	X	÷	
Notes:	MS MSD	phlecker				
Weather:	Temperature <sup>o</sup> F_	205	Sunny Cloudy	Rain Snow		
Revised 08/15/07						

Client/Site: Monitoring Location:	Quanta Resources			Project No.: Date:	2015007
Source Description:	MW-10			Sampler: MT	
Well & Water Level I	Data:		l Depth of Well: Depth to Water: Column (LWC):	34.60feet <u>24.69</u> feet 9.90 feet	
Purge Volume Calcula Well Diameter (i 1		LWC * 0.041	2	Gallons	
1.25 1.5 2 3 4 6		LWC * 0.064 LWC * 0.092 LWC * 0.163 LWC * 0.367 LWC * 0.653 LWC * 1.469	*3 = *3 = *3 = *3 =	Gallons Gallons Gallons Gallons Gallons Gallons	
Free Product Check:	Free Pı Measured Thickı	roduct Present: ness/Comment: _	Yes	No no bailer	
Purge Data:	Purge Date: _	12/5/18		e -	
	<b>Purging Time:</b>	From:	#1:5D	То:	2-40
7	Гуре of Purging Eq Purged Wa	uipment Used: _ ter Comments: _	Bladder Cleur -	Pump some oder - pai	tu
Sampling Data:	Depth to Wate	er at Sampling: _	24.50	fee	t
	Color of Sample: Turbidity:	slight	Sample Date: Sample Time:	110:40	>
	ype of Sampling Eq Present During Sam		Bludeb r Odor	Рл <i>р</i>	
Notes:			Sheen Free Product None	<u>×</u>	
Weather:	Temperature <sup>o</sup> F	10 (	Sunny Cloudy	Rain Snow	

Client/Site:	Quanta Resources			<b>Project No.:</b>	
<b>Monitoring Location:</b>				Date:	12/5/2018
Source Description:	MW-12			Sampler:	MTM/DTH
Well & Water Level Data:		Tota	l Depth of Well:	37.42	feet
		Initial	Depth to Water:	26.80	feet
	Le	ength of Water (	Column (LWC):		feet
Purge Volume Calcula	ation:				
Well Diameter (in		Calculated W	ell Volume To Be	e Removed	
1	,	LWC * 0.041		Gallons	
1.25			* 3 =	Gallons	
1.5			* 3 =	Gallons	
2		LWC * 0.163 * 3 =		Gallons	
3		LWC $* 0.367 * 3 =$		Gallons	
4		LWC $* 0.653 * 3 =$		Gallons	
6		LWC * 1.469	* 3 =	Gallons	
Free Product Check:	Free Pr	oduct Present:	Yes	No	
	Measured Thick			C	
Purge Data:	Purge Date:	12/5/18			
C	Purging Time:	1.4.35	12:00	То•	1:20
	• •	-			1. 20
ĩ	ype of Purging Eq Purged Wa	uipment Used:	Blackder	Punp	
	Purged Wa	ter Comments:	Cleare	-	
Sampling Data:	Depth to Wate	er at Sampling:			feet
(	Color of Sample:	cluc	Sample Date:	12/5/18	
,		slight	Sample Time:	14:50	
		Slight			
Ту	pe of Sampling Eq	uipment Used:	Bluiller	Pung	
<b>Field Indicators P</b>	resent During Sam	ple Collection:	Odor		
			Sheen		
			Free Product		
			None	×	
Notes:					
÷					
Weather:	Temperature <sup>o</sup> F	70	Sunny Cloudy	Rain Snow	
		_	No. and the second seco		

# LABORATORY REPORT



### Dayton, NJ

The results set forth herein are provided by SGS North America Inc.

e-Hardcopy 2.0 Automated Report

12/26/18

# **Technical Report for**

### **Plumley Environmental Engineers**

Quanta Resources, Lodi Street, Syracuse, NY

2015127.006

SGS Job Number: JC79316



Sampling Date: 12/05/18

Report to:

**Plumley Environmental Engineers** 

dhudson@plumleyeng.com

ATTN: Derk Hudson

### Total number of pages in report: 50



MATT

Test results contained within this data package meet the requirements of the National Environmental Laboratory Accreditation Program and/or state specific certification programs as applicable.

Brian McGuire General Manager

Client Service contact: Thelma Flaherty 732-329-0200

Certifications: NJ(12129), NY(10983), CA, CT, FL, IL, IN, KS, KY, LA, MA, MD, ME, MN, NC, OH VAP (CL0056), AK (UST-103), AZ (AZ0786), PA, RI, SC, TX, UT, VA, WV, DoD ELAP (ANAB L2248)

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Please share your ideas about how we can serve you better at: EHS.US.CustomerCare@sgs.com

1 of 50

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# **Sample Summary**

Plumley Environmental Engineers

Job No: JC79316

Quanta Resources, Lodi Street, Syracuse, NY Project No: 2015127.006

Sample Number	Collected Date	Time By	Received	Matr Code		Client Sample ID
JC79316-1	12/05/18	12:30 MM	12/07/18	AQ	Ground Water	MW-9
JC79316-1D	12/05/18	12:30 MM	12/07/18	AQ	Water Dup/MSD	MW-9 MSD
JC79316-1S	12/05/18	12:30 MM	12/07/18	AQ	Water Matrix Spike	MW-9 MS
JC79316-2	12/05/18	12:50 MM	12/07/18	AQ	Equipment Blank	EQUIPMENT BLANK
JC79316-3	12/05/18	13:01 MM	12/07/18	AQ	Ground Water	MW-6
JC79316-4	12/05/18	13:25 MM	12/07/18	AQ	Ground Water	MW-5
JC79316-5	12/05/18	13:47 MM	12/07/18	AQ	Ground Water	MW-2
JC79316-6	12/05/18	14:00 MM	12/07/18	AQ	Ground Water	MW-1-D
JC79316-7	12/05/18	14:50 MM	12/07/18	AQ	Ground Water	MW-12
JC79316-8	12/05/18	16:40 MM	12/07/18	AQ	Ground Water	MW-10





## **CASE NARRATIVE / CONFORMANCE SUMMARY**

Client:	Plumley Environmental Engineers	Job No	JC79316
Site:	Quanta Resources, Lodi Street, Syracuse, NY	Report Date	12/26/2018 2:40:17 P

On 12/07/2018, 8 Sample(s), 0 Trip Blank(s) and 0 Field Blank(s) were received at SGS North America Inc. at a maximum corrected temperature of 1.8 C. Samples were intact and chemically preserved, unless noted below. A SGS North America Inc. Job Number of JC79316 was assigned to the project. Laboratory sample ID, client sample ID and dates of sample collection are detailed in the report's Results Summary Section.

Specified quality control criteria were achieved for this job except as noted below. For more information, please refer to the analytical results and QC summary pages.

Compounds qualified as out of range in the continuing calibration summary report are acceptable as per method requirements when there is a high bias but the sample result is non-detect.

#### MS Volatiles By Method SW846 8260C

Matrix: AQ	Batch ID: V2V2263	
		,

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JC79316-1MS, JC79316-1MSD were used as the QC samples indicated.
- Blank Spike Recovery(s) for 1,1,2,2-Tetrachloroethane are outside control limits. High percent recoveries and no associated positive reported in the QC batch.
- RPD(s) for MSD for Bromomethane are outside control limits for sample JC79316-1MSD. Outside control limits due to matrix interference.
- JC79316-6 for Dichlorodifluoromethane: Associated CCV outside of control limits low.
- JC79316-4 for 1,1,2,2-Tetrachloroethane: This compound in BS is outside in house QC limits bias high.
- JC79316-1 for Bromomethane: Associated CCV outside of control limits low.
- JC79316-3 for Bromomethane: Associated CCV outside of control limits low.
- JC79316-3 for Dichlorodifluoromethane: Associated CCV outside of control limits low.
- JC79316-3 for 1,1,2,2-Tetrachloroethane: This compound in BS is outside in house QC limits bias high.
- JC79316-4 for Bromomethane: Associated CCV outside of control limits low.
- JC79316-6 for 1,1,2,2-Tetrachloroethane: This compound in BS is outside in house QC limits bias high.
- JC79316-6 for Bromomethane: Associated CCV outside of control limits low.
- JC79316-1 for 1,1,2,2-Tetrachloroethane: This compound in BS is outside in house QC limits bias high.
- JC79316-7 for Dichlorodifluoromethane: Associated CCV outside of control limits low.
- JC79316-7 for 1,1,2,2-Tetrachloroethane: This compound in BS is outside in house QC limits bias high.
- JC79316-4 for Dichlorodifluoromethane: Associated CCV outside of control limits low.
- JC79316-8 for Dichlorodifluoromethane: Associated CCV outside of control limits low.
- JC79316-8 for 1,1,2,2-Tetrachloroethane: This compound in BS is outside in house QC limits bias high.
- JC79316-7 for Bromomethane: Associated CCV outside of control limits low.
- JC79316-1 for Dichlorodifluoromethane: Associated CCV outside of control limits low.
- JC79316-8 for Bromomethane: Associated CCV outside of control limits low.

#### Matrix: AQ Batch ID: V2V2267

- All samples were analyzed within the recommended method holding time.
- Sample(s) JC79319-15MS, JC79319-15MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.
- Blank Spike Recovery(s) for 1,1,2,2-Tetrachloroethane are outside control limits. High percent recoveries and no associated positive reported in the QC batch.

#### Wednesday, December 26, 2018

#### MS Volatiles By Method SW846 8260C

Matrix: AQ	Batch ID: V2V2267	
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- Matrix Spike Recovery(s) for 1,1-Dichloroethane are outside control limits. Outside control limits due to high level in sample relative to spike amount.
- RPD(s) for MSD for Bromomethane are outside control limits for sample JC79319-15MSD. Probable cause due to sample homogeneity.
- JC79316-5 for 1,1,2,2-Tetrachloroethane: Associated CCV outside of control limits high, sample was ND. This compound in BS is outside in house QC limits bias high.
- JC79319-15MSD for Bromomethane: Outside control limits due to matrix interference.
- JC79316-5 for Bromomethane: Associated CCV outside of control limits low.
- JC79316-5 for 1,2-Dichloroethane: Associated CCV outside of control limits high, sample was ND.
- V2V2267-BS for 1,1,2,2-Tetrachloroethane: High percent recoveries and no associated positive reported in the QC batch.

#### MS Semi-volatiles By Method EPA 537M BY ID

	Matrix: AQ	Batch ID: F:OP73036	
- 27	THE LASS EDA 527M DV ID		

- The data for EPA 537M BY ID meets quality control requirements.
- JC79316-1: Dilution required due to matrix interference. Analysis performed at SGS Orlando, FL.
- JC79316-2: Analysis performed at SGS Orlando, FL.
- JC79316-8: Dilution required due to matrix interference. Analysis performed at SGS Orlando, FL.
- JC79316-7: Analysis performed at SGS Orlando, FL.
- JC79316-1 for Perfluorooctanoic acid: Associated CCV outside of control limits high.
- JC79316-8 for Perfluorooctanoic acid: Associated CCV outside of control limits high, sample was ND.

#### MS Semi-volatiles By Method SW846 8270D BY SIM

	Matrix: AQ	Batch ID: OP17214A	
. 1			

All samples were extracted within the recommended method holding time.

All method blanks for this batch meet method specific criteria.

#### GC/LC Semi-volatiles By Method SW846 8082A

	Matrix: AQ	Batch ID:	OP17274
_	A11 1 1		11 11

- All samples were extracted within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JC79316-3 have surrogates outside control limits. Probable cause due to matrix interference.
- JC79316-5 for Aroclor 1248: More than 40 % RPD for detected concentrations between the two GC columns.
- JC79316-3 for Decachlorobiphenyl: Outside the QC limits. There is no sample left to re-extract.

SGS North America Inc. certifies that data reported for samples received, listed on the associated custody chain or analytical task order, were produced to specifications meeting the Quality System precision, accuracy and completeness objectives except as noted.

Estimated non-standard method measurement uncertainty data is available on request, based on quality control bias and implicit for standard methods. Acceptable uncertainty requires tested parameter quality control data to meet method criteria.

SGS North America Inc. is not responsible for data quality assumptions if partial reports are used and recommends that this report be used in its entirety. Data release is authorized by SGS North America Inc indicated via signature on the report cover

## SAMPLE DELIVERY GROUP CASE NARRATIVE

Client:	SGS Dayton, NJ	Job No	JC79316
Site:	PLUMNYB: Quanta Resources, Lodi Street, Syracuse, NY	Report Date	12/26/2018 1:51:04

4 Samples were collected on 12/05/2018 and were received at SGS North America Inc - Orlando on 12/07/2018 properly preserved, at 3.8 Deg. C and intact. These samples received an SGS Orlando job number of JC79316. A listing of the Laboratory Sample ID, Client Sample ID and dates of collection are presented in the Results Summary Section. Except as noted below, all method specified calibrations and quality control performance criteria were met for this job. For more information, please refer to QC summary pages.

#### MS Semi-Volatiles By Method EPA 537M BY ID

#### Matrix: AQ Batch ID: OP73036

All samples were extracted within the recommended method holding time.

All samples were analyzed within the recommended method holding time.

Sample(s) JC79316-1MS, JC79316-1MSD were used as the QC samples indicated.

All method blanks for this batch meet method specific criteria.

JC79316-1 for Perfluorooctanoic acid: Associated CCV outside of control limits high.

JC79316-1: Dilution required due to matrix interference.

JC79316-8 for Perfluorooctanoic acid: Associated CCV outside of control limits high, sample was ND.

JC79316-8: Dilution required due to matrix interference.

SGS Orlando certifies that this report meets the project requirements for analytical data produced for the samples as received at SGS Orlando and as stated on the COC. SGS Orlando certifies that the data meets the Data Quality Objectives for precision, accuracy and completeness as specified in the SGS Orlando Quality Manual except as noted above. This report is to be used in its entirety. SGS Orlando is not responsible for any assumptions of data quality if partial data packages are used.

Narrative prepared by:

Ariel Hartney, Client Services (Signature on File)

# **Summary of Hits**

Job Number:	JC79316
Account:	Plumley Environmental Engineers
Project:	Quanta Resources, Lodi Street, Syracuse, NY
Collected:	12/05/18

Lab Sample ID Client Sample ID Analyte	Result/ Qual	RL	MDL	Units	Method
JC79316-1 MW-9					
Benzene	0.66	0.50	0.43	ug/l	SW846 8260C
Chlorobenzene	5.8	1.0	0.56	ug/l	SW846 8260C
1,2-Dichlorobenzene	1.0	1.0	0.53	ug/l	SW846 8260C
1,3-Dichlorobenzene	0.62 J	1.0	0.54	ug/l	SW846 8260C
1,4-Dichlorobenzene	1.6	1.0	0.51	ug/l	SW846 8260C
Isopropylbenzene	0.72 J	1.0	0.65	ug/l	SW846 8260C
Toluene	0.66 J	1.0	0.53	ug/l	SW846 8260C
Perfluorobutanoic acid <sup>a</sup>	10.5 J	15	3.8	ng/l	EPA 537M BY ID
Perfluorohexanoic acid <sup>a</sup>	7.03 J	7.7	1.9	ng/l	EPA 537M BY ID
Perfluoroheptanoic acid <sup>a</sup>	2.09 J	3.8	1.9	ng/l	EPA 537M BY ID
Perfluorooctanoic acid b	23.6	3.8	1.9	ng/l	EPA 537M BY ID
Perfluorobutanesulfonic acid <sup>a</sup>	1.93 J	3.8	1.9	ng/l	EPA 537M BY ID
Perfluorohexanesulfonic acid <sup>a</sup>	2.87 J	3.8	1.9	ng/l	EPA 537M BY ID
Perfluorooctanesulfonic acid <sup>a</sup>	3.40 J	3.8	2.9	ng/l	EPA 537M BY ID
1,4-Dioxane	2.34	0.095	0.046	ug/l	SW846 8270D BY SIM

#### JC79316-2 EQUIPMENT BLANK

No hits reported in this sample.

#### JC79316-3 MW-6

Benzene	1.7	0.50	0.43	ug/l	SW846 8260C
Chlorobenzene	23.9	1.0	0.56	ug/l	SW846 8260C
1,2-Dichlorobenzene	4.5	1.0	0.53	ug/l	SW846 8260C
1,3-Dichlorobenzene	2.4	1.0	0.54	ug/l	SW846 8260C
1,4-Dichlorobenzene	2.7	1.0	0.51	ug/l	SW846 8260C
1,1-Dichloroethane	0.61 J	1.0	0.57	ug/l	SW846 8260C
JC79316-4 MW-5					
Bromodichloromethane	0.82 J	1.0	0.58	ug/l	SW846 8260C
Chloroform	17.4	1.0	0.50	ug/l	SW846 8260C
				0	
JC79316-5 MW-2					
Benzene	1.4	0.50	0.43	ug/l	SW846 8260C
Chlorobenzene	0.97 J	1.0	0.56	ug/l	SW846 8260C
Chloroethane	0.89 J	1.0	0.73	ug/l	SW846 8260C
1,2-Dichlorobenzene	3.1	1.0	0.53	ug/l	SW846 8260C
1,3-Dichlorobenzene	2.5	1.0	0.54	ug/l	SW846 8260C
1,4-Dichlorobenzene	5.9	1.0	0.51	ug/l	SW846 8260C
1,1-Dichloroethane	0.73 J	1.0	0.57	ug/l	SW846 8260C



# **Summary of Hits**

Job Number:	JC79316
Account:	Plumley Environmental Engineers
Project:	Quanta Resources, Lodi Street, Syracuse, NY
Collected:	12/05/18

Lab Sample ID Client Sample ID Analyte	Result/ Qual	RL	MDL	Units	Method
1,2-Dichloroethene (total)	1.0	1.0	0.51	ug/l	SW846 8260C
Trichloroethene	1.2	1.0	0.53	ug/l	SW846 8260C
Aroclor 1248 <sup>c</sup>	0.75	0.24	0.061	ug/l	SW846 8082A
Aroclor 1254	1.1	0.24	0.20	ug/l	SW846 8082A
Aroclor 1260	1.2	0.24	0.073	ug/l	SW846 8082A
JC79316-6 MW-1-D					
Benzene	0.52	0.50	0.43	ug/l	SW846 8260C
Chlorobenzene	0.57 J	1.0	0.56	ug/l	SW846 8260C
1,2-Dichlorobenzene	0.59 J	1.0	0.53	ug/l	SW846 8260C
1,1-Dichloroethane	2.3	1.0	0.57	ug/l	SW846 8260C
1,2-Dichloroethene (total)	1.5	1.0	0.51	ug/l	SW846 8260C
Vinyl chloride	24.0	1.0	0.79	ug/l	SW846 8260C
JC79316-7 MW-12					
Benzene	0.46 J	0.50	0.43	ug/l	SW846 8260C
sec-Butylbenzene	1.4 J	2.0	0.62	ug/l	SW846 8260C
tert-Butylbenzene	0.96 J	2.0	0.69	ug/l	SW846 8260C
Chlorobenzene	11.2	1.0	0.56	ug/l	SW846 8260C
1,2-Dichlorobenzene	3.4	1.0	0.53	ug/l	SW846 8260C
1,4-Dichlorobenzene	1.8	1.0	0.51	ug/l	SW846 8260C
Isopropylbenzene	3.9	1.0	0.65	ug/l	SW846 8260C
n-Propylbenzene	1.3 J	2.0	0.60	ug/l	SW846 8260C
Perfluorobutanoic acid d	4.87 J	8.0	2.0	ng/l	EPA 537M BY ID
Perfluorohexanoic acid <sup>d</sup>	1.31 J	4.0	1.0	ng/l	EPA 537M BY ID
Perfluorooctanoic acid <sup>d</sup>	3.41	2.0	1.0	ng/l	EPA 537M BY ID
Perfluorooctanesulfonic acid <sup>d</sup>	6.12	2.0	1.5	ng/l	EPA 537M BY ID
1,4-Dioxane	0.239	0.095	0.046	ug/l	SW846 8270D BY SIM
JC79316-8 MW-10					
Benzene	2.3	0.50	0.43	ug/l	SW846 8260C
n-Butylbenzene	0.63 J	2.0	0.52	ug/l	SW846 8260C
sec-Butylbenzene	1.3 J	2.0	0.62	ug/l	SW846 8260C
1,2-Dichlorobenzene	2.6	1.0	0.53	ug/l	SW846 8260C
1,4-Dichlorobenzene	1.1	1.0	0.51	ug/l	SW846 8260C
1,2-Dichloroethene (total)	0.93 J	1.0	0.51	ug/l	SW846 8260C
Isopropylbenzene	2.3	1.0	0.65	ug/l	SW846 8260C
n-Propylbenzene	1.9 J	2.0	0.60	ug/l	SW846 8260C
1,4-Dioxane	0.225	0.10	0.00	ug/l	SW846 8270D BY SIM
Aroclor 1254	0.225	0.10	0.20	ug/l	SW846 8082A
	0.20	0.21	0.20	~B' 1	5010 000 <b>2</b> 11

(a) Dilution required due to matrix interference. Analysis performed at SGS Orlando, FL.

# **Summary of Hits**

Job Number:	JC79316
Account:	Plumley Environmental Engineers
Project:	Quanta Resources, Lodi Street, Syracuse, NY
Collected:	12/05/18

Lab Sample ID	<b>Client Sample ID</b>	Result/					
Analyte		Qual	RL	MDL	Units	Method	

(b) Dilution required due to matrix interference. Analysis performed at SGS Orlando, FL. Associated CCV outside of control limits high.

(c) More than 40 % RPD for detected concentrations between the two GC columns.

(d) Analysis performed at SGS Orlando, FL.

ω



Dayton, NJ

Section 4

Sample Results

Report of Analysis



4



Client Sa Lab Sam Matrix: Method: Project:	ple ID: JC A SV	W-9 279316-1 Q - Ground W W846 8260C uanta Resource	ater es, Lodi Street, Syrad	cuse, N		Date Sampled: Date Received: Percent Solids:	/ -/
Run #1 Run #2	<b>File ID</b> 2V55851.1	<b>DF</b> D 1	<b>Analyzed</b> 12/12/18 09:16	By JP	<b>Prep Date</b> n/a	<b>Prep Batc</b> n/a	h Analytical Batch V2V2263
Run #1 Run #2	Purge Vol 5.0 ml	ume					

#### VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	6.0	ug/l	
71-43-2	Benzene	0.66	0.50	0.43	ug/l	
108-86-1	Bromobenzene	ND	1.0	0.55	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.58	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>a</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	6.9	ug/l	
104-51-8	n-Butylbenzene	ND	2.0	0.52	ug/l	
135-98-8	sec-Butylbenzene	ND	2.0	0.62	ug/l	
98-06-6	tert-Butylbenzene	ND	2.0	0.69	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.95	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	5.8	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	ND	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
95-49-8	o-Chlorotoluene	ND	2.0	0.63	ug/l	
106-43-4	p-Chlorotoluene	ND	2.0	0.60	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	1.2	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	1.0	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	0.62	1.0	0.54	ug/l	J
106-46-7	1,4-Dichlorobenzene	1.6	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane <sup>a</sup>	ND	2.0	1.4	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
540-59-0	1,2-Dichloroethene (total)	ND	1.0	0.51	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
142-28-9	1,3-Dichloropropane	ND	1.0	0.43	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



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JC79316

E = Indicates value exceeds calibration range

Client Sample ID:	MW-9		
Lab Sample ID:	JC79316-1	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	SW846 8260C	Percent Solids:	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

**VOA Special List** 

CAS No.	Compound	Result	RL	MDL	Units	Q
594-20-7	2,2-Dichloropropane	ND	1.0	0.52	ug/l	
563-58-6	1,1-Dichloropropene	ND	1.0	0.82	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
87-68-3	Hexachlorobutadiene	ND	2.0	0.56	ug/l	
98-82-8	Isopropylbenzene	0.72	1.0	0.65	ug/l	J
99-87-6	p-Isopropyltoluene	ND	2.0	0.66	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
74-95-3	Methylene bromide	ND	1.0	0.48	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
91-20-3	Naphthalene	ND	5.0	0.98	ug/l	
103-65-1	n-Propylbenzene	ND	2.0	0.60	ug/l	
100-42-5	Styrene	ND	1.0	0.70	ug/l	
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0	0.60	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane <sup>b</sup>	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.90	ug/l	
108-88-3	Toluene	0.66	1.0	0.53	ug/l	J
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.84	ug/l	
96-18-4	1,2,3-Trichloropropane	ND	2.0	0.70	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	2.0	1.0	ug/l	
108-67-8	1,3,5-Trimethylbenzene	ND	2.0	1.0	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.79	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Lim	its	
1868-53-7	Dibromofluoromethane	98%		80-1	20%	
17060-07-0	1,2-Dichloroethane-D4	109%		81-1	24%	
2037-26-5	Toluene-D8	98%		80-1	20%	
460-00-4	4-Bromofluorobenzene	101%		80-1	20%	

(a) Associated CCV outside of control limits low.

(b) This compound in BS is outside in house QC limits bias high.

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound

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E = Indicates value exceeds calibration range

			Report	of Ana	alysis			Page 1 of 1
Client Sam Lab Sampl Matrix: Method: Project:	e ID: JC793 AQ - SW84	16-1 Ground Wate 6 8270D BY	r SIM SW846 35 Lodi Street, Syra			Date	L .	2/05/18 2/07/18 ′a
Run #1 Run #2	<b>File ID</b> 3P73448.D	<b>DF</b> 1	<b>Analyzed</b> 12/13/18 02:49	By SA	<b>Prep D</b> 12/11/1		<b>Prep Batch</b> OP17214A	<b>Analytical Batch</b> E3P3464
Run #1 Run #2	<b>Initial Volum</b> 1050 ml	e Final Vol 1.0 ml	lume					
CAS No.	Compound		Result	RL	MDL	Units	Q	
123-91-1	1,4-Dioxane		2.34	0.095	0.046	ug/l		
CAS No.	Surrogate Re	ecoveries	Run# 1	Run# 2	Lim	its		
4165-60-0 321-60-8 1718-51-0	Nitrobenzene 2-Fluorobiph Terphenyl-d1	enyl	56% 45% 46%		29-1 23-1 22-1	22%		

ND = Not detected MDL = Method Detection Limit

- RL = Reporting Limit
- E = Indicates value exceeds calibration range
- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

	nple ID: MW-9	<i>c</i> 1					2/05/10
Lab Samp	ole ID: JC7931	6-1			Date	Sampled: 1	2/05/18
Matrix:	AQ - G	round Wa	iter		Date	Received: 1	2/07/18
Method:	EPA 53	37M BY I	D EPA 537 MOD		Perc	ent Solids: n	n/a
Project:	Quanta	Resource	s, Lodi Street, Syrad	cuse, N	Y		
	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup> Run #2	<b>File ID</b> 2Q25412.D	<b>DF</b> 2	Analyzed 12/21/18 20:05	•	<b>Prep Date</b> 12/14/18 08:45	-	<b>Analytical Batch</b> F:S2Q394

Initial Volume Final Vol Run #1 260 ml 1.0 ml Run #2

#### PFAS List

CAS No.	Compound	Result	RL	MDL	Units	Q
375-22-4	Perfluorobutanoic acid	10.5	15	3.8	ng/l	J
2706-90-3	Perfluoropentanoic acid	ND	7.7	2.9	ng/l	
307-24-4	Perfluorohexanoic acid	7.03	7.7	1.9	ng/l	J
375-85-9	Perfluoroheptanoic acid	2.09	3.8	1.9	ng/l	J
335-67-1	Perfluorooctanoic acid b	23.6	3.8	1.9	ng/l	
375-95-1	Perfluorononanoic acid	ND	3.8	1.9	ng/l	
335-76-2	Perfluorodecanoic acid	ND	7.7	1.9	ng/l	
2058-94-8	Perfluoroundecanoic acid	ND	7.7	1.9	ng/l	
307-55-1	Perfluorododecanoic acid	ND	7.7	2.9	ng/l	
72629-94-8	Perfluorotridecanoic acid	ND	7.7	1.9	ng/l	
376-06-7	Perfluorotetradecanoic acid	ND	7.7	1.9	ng/l	
375-73-5	Perfluorobutanesulfonic acid	1.93	3.8	1.9	ng/l	J
355-46-4	Perfluorohexanesulfonic acid	2.87	3.8	1.9	ng/l	J
375-92-8	Perfluoroheptanesulfonic acid	ND	7.7	1.9	ng/l	
1763-23-1	Perfluorooctanesulfonic acid	3.40	3.8	2.9	ng/l	J
335-77-3	Perfluorodecanesulfonic acid	ND	7.7	1.9	ng/l	
754-91-6	PFOSA	ND	7.7	1.9	ng/l	
2355-31-9	MeFOSAA	ND	38	7.7	ng/l	
2991-50-6	EtFOSAA	ND	38	7.7	ng/l	
27619-97-2	6:2 Fluorotelomer sulfonate	ND	15	3.8	ng/l	
39108-34-4	8:2 Fluorotelomer sulfonate	ND	15	3.8	ng/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Lim	its	
	13C4-PFBA	79%		30-1	40%	
	13C5-PFPeA	90%		40-1	40%	
	13C5-PFHxA	94%		50-1	50%	
	13C4-PFHpA	103%		50-1	50%	
	13C8-PFOA	120%		50-1	50%	
	13C9-PFNA	112%		50-1	50%	
	13C6-PFDA	97%		50-1	50%	
	13C7-PFUnDA	99%		50-1	50%	

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound

Page 1 of 2

Client Sample ID:	MW 0		
Chefft Sample ID.	IVI VV -9		
Lab Sample ID:	JC79316-1	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	EPA 537M BY ID EPA 537 MOD	<b>Percent Solids:</b>	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

#### **PFAS List**

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
	13C2-PFDoDA	108%		50-150%
	13C2-PFTeDA	123%		40-150%
	13C3-PFBS	94%		50-150%
	13C3-PFHxS	99%		50-150%
	13C8-PFOS	101%		50-150%
	13C8-FOSA	50%		30-140%
	d3-MeFOSAA	124%		50-150%
	13C2-6:2FTS	148%		50-150%
	13C2-8:2FTS	112%		50-150%

(a) Dilution required due to matrix interference. Analysis performed at SGS Orlando, FL.

(b) Associated CCV outside of control limits high.

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound

Page 2 of 2



			Report	of An	alysis			Page 1 of 1
Client Sam Lab Samp Matrix: Method: Project:	le ID: JC79310 AQ - Gi SW846	ound Water 8082A SW		cuse, NY	7	Date	Received:	12/05/18 12/07/18 n/a
Run #1 Run #2	<b>File ID</b> XX240797.D	<b>DF</b> 1	<b>Analyzed</b> 12/13/18 23:09	<b>By</b> CP	<b>Prep Date</b> 12/13/18		Prep Batch OP17274	Analytical Batch GXX6552
Run #1 Run #2	<b>Initial Volume</b> 1040 ml	<b>Final Volu</b> 5.0 ml	ime					
PCB List								
CAS No.	Compound		Result	RL	MDL	Units	Q	

CAB III.	Compound	KtSuit	<b>KL</b>	MDL	Omts	V
12674-11-2	Aroclor 1016	ND	0.24	0.094	ug/l	
11104-28-2	Aroclor 1221	ND	0.24	0.20	ug/l	
11141-16-5	Aroclor 1232	ND	0.24	0.12	ug/l	
53469-21-9	Aroclor 1242	ND	0.24	0.11	ug/l	
12672-29-6	Aroclor 1248	ND	0.24	0.061	ug/l	
11097-69-1	Aroclor 1254	ND	0.24	0.20	ug/l	
11096-82-5	Aroclor 1260	ND	0.24	0.073	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limit	ts	
877-09-8	Tetrachloro-m-xylene	40%		11-16	66%	
877-09-8	Tetrachloro-m-xylene	35%		11-16	6%	
2051-24-3	Decachlorobiphenyl	17%		10-15	50%	
2051-24-3	Decachlorobiphenyl	19%		10-15	50%	
	1 2					

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





Project:	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 <sup>a</sup> Run #2	3Q558.D	1	12/18/18 22:30	AFL	12/14/18 08:45	F:OP73036	F:S3Q9

Run #1 260 ml 1.

Run #2

1.0 ml

#### **PFAS List**

CAS No.	Compound	Result	RL	MDL	Units	Q
375-22-4	Perfluorobutanoic acid	ND	7.7	1.9	ng/l	
2706-90-3	Perfluoropentanoic acid	ND	3.8	1.4	ng/l	
307-24-4	Perfluorohexanoic acid	ND	3.8	0.96	ng/l	
375-85-9	Perfluoroheptanoic acid	ND	1.9			
335-67-1	Perfluorooctanoic acid	ND	1.9	0.96	ng/l	
375-95-1	Perfluorononanoic acid	ND	1.9	0.96	ng/l	
335-76-2	Perfluorodecanoic acid	ND	3.8	0.96	ng/l	
2058-94-8	Perfluoroundecanoic acid	ND	3.8	0.96	ng/l	
307-55-1	Perfluorododecanoic acid	ND	3.8	1.4	ng/l	
72629-94-8	Perfluorotridecanoic acid	ND	3.8	0.96	ng/l	
376-06-7	Perfluorotetradecanoic acid	ND	3.8	0.96	ng/l	
375-73-5	Perfluorobutanesulfonic acid	ND	1.9	0.96	ng/l	
355-46-4	Perfluorohexanesulfonic acid	ND	1.9	0.96	ng/l	
375-92-8	Perfluoroheptanesulfonic acid	ND	3.8	0.96	ng/l	
1763-23-1	Perfluorooctanesulfonic acid	ND	1.9	1.4	ng/l	
335-77-3	Perfluorodecanesulfonic acid	ND	3.8	0.96	ng/l	
754-91-6	PFOSA	ND	3.8	0.96	ng/l	
2355-31-9	MeFOSAA	ND	19	3.8	ng/l	
2991-50-6	EtFOSAA	ND	19	3.8	ng/l	
27619-97-2	6:2 Fluorotelomer sulfonate	ND	7.7	1.9	ng/l	
39108-34-4	8:2 Fluorotelomer sulfonate	ND	7.7	1.9	ng/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Lim	its	
	13C4-PFBA	70%		30-1	40%	
	13C5-PFPeA	72%		40-1	40%	
	13C5-PFHxA	77%		50-1	50%	
	13C4-PFHpA	80%		50-1	50%	
	13C8-PFOA	84%		50-1	50%	
	13C9-PFNA	86%		50-1	50%	
	13C6-PFDA	103%		50-1	50%	
	13C7-PFUnDA	89%		50-1	50%	

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



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Client Sample ID:	EQUIPMENT BLANK		
Lab Sample ID:	JC79316-2	Date Sampled:	12/05/18
Matrix:	AQ - Equipment Blank	Date Received:	12/07/18
Method:	EPA 537M BY ID EPA 537 MOD	Percent Solids:	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

#### **PFAS List**

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CAS No.	Surrogate Recoveries	<b>Run#</b> 1	Run# 2	Limits
	13C2-PFDoDA	76%		50-150%
	13C2-PFTeDA	72%		40-150%
	13C3-PFBS	72%		50-150%
	13C3-PFHxS	76%		50-150%
	13C8-PFOS	77%		50-150%
	13C8-FOSA	92%		30-140%
	d3-MeFOSAA	92%		50-150%
	13C2-6:2FTS	82%		50-150%
	13C2-8:2FTS	96%		50-150%

(a) Analysis performed at SGS Orlando, FL.

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound

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Client San Lab Samj Matrix: Method: Project:	ple ID: JC79 AQ - SW8	316-3 Ground Wa 46 8260C	ater es, Lodi Street, Syrad	cuse, N	I P	Date Sampled: 1 Date Received: 1 Percent Solids: n	
Run #1 Run #2	<b>File ID</b> 2V55852.D	<b>DF</b> 1	<b>Analyzed</b> 12/12/18 09:41	<b>By</b> JP	<b>Prep Date</b> n/a	<b>Prep Batch</b> n/a	Analytical Batch V2V2263
Run #1 Run #2	<b>Purge Volum</b> 5.0 ml	ie					

#### **VOA Special List**

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	6.0	ug/l	
71-43-2	Benzene	1.7	0.50	0.43	ug/l	
108-86-1	Bromobenzene	ND	1.0	0.55	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.58	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>a</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	6.9	ug/l	
104-51-8	n-Butylbenzene	ND	2.0	0.52	ug/l	
135-98-8	sec-Butylbenzene	ND	2.0	0.62	ug/l	
98-06-6	tert-Butylbenzene	ND	2.0	0.69	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.95	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	23.9	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	ND	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
95-49-8	o-Chlorotoluene	ND	2.0	0.63	ug/l	
106-43-4	p-Chlorotoluene	ND	2.0	0.60	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	1.2	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	4.5	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	2.4	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	2.7	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane <sup>a</sup>	ND	2.0	1.4	ug/l	
75-34-3	1,1-Dichloroethane	0.61	1.0	0.57	ug/l	J
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
540-59-0	1,2-Dichloroethene (total)	ND	1.0	0.51	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
142-28-9	1,3-Dichloropropane	ND	1.0	0.43	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound

Page 1 of 2



E = Indicates value exceeds calibration range

J = Indicates an estimated value

Client Sample ID:	MW-6		
Lab Sample ID:	JC79316-3	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	SW846 8260C	Percent Solids:	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

**VOA Special List** 

CAS No.	Compound	Result	RL	MDL	Units	Q	
594-20-7	2,2-Dichloropropane	ND	1.0	0.52	ug/l		
563-58-6	1,1-Dichloropropene	ND	1.0	0.82	ug/l		
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l		
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l		
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l		
87-68-3	Hexachlorobutadiene	ND	2.0	0.56	ug/l		
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l		
99-87-6	p-Isopropyltoluene	ND	2.0	0.66	ug/l		
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l		
74-95-3	Methylene bromide	ND	1.0	0.48	ug/l		
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l		
91-20-3	Naphthalene	ND	5.0	0.98	ug/l		
103-65-1	n-Propylbenzene	ND	2.0	0.60	ug/l		
100-42-5	Styrene	ND	1.0	0.70	ug/l		
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0	0.60	ug/l		
79-34-5	1,1,2,2-Tetrachloroethane <sup>b</sup>	ND	1.0	0.65	ug/l		
127-18-4	Tetrachloroethene	ND	1.0	0.90	ug/l		
108-88-3	Toluene	ND	1.0	0.53	ug/l		
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l		
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l		
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l		
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l		
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l		
75-69-4	Trichlorofluoromethane	ND	2.0	0.84	ug/l		
96-18-4	1,2,3-Trichloropropane	ND	2.0	0.70	ug/l		
95-63-6	1,2,4-Trimethylbenzene	ND	2.0	1.0	ug/l		
108-67-8	1,3,5-Trimethylbenzene	ND	2.0	1.0	ug/l		
75-01-4	Vinyl chloride	ND	1.0	0.79	ug/l		
	m,p-Xylene	ND	1.0	0.78	ug/l		
95-47-6	o-Xylene	ND	1.0	0.59	ug/l		
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Lim	its		
1868-53-7	Dibromofluoromethane	98%		80-1	20%		
17060-07-0	1,2-Dichloroethane-D4	107%		81-1	24%		
2037-26-5	Toluene-D8	98%		80-1	20%		
460-00-4				80-120%			

(a) Associated CCV outside of control limits low.

(b) This compound in BS is outside in house QC limits bias high.

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound

Page 2 of 2



E = Indicates value exceeds calibration range

				Repor	rt of Ana	alysis			Page 1 of 1
Client Sample ID: Lab Sample ID: Matrix: Method: Project:		SW846	ound Wate 8082A SV	r V846 3510C Lodi Street, Sy	racuse, NY		Date	Received: 1	2/05/18 2/07/18 /a
Run #1 Run #2	File ID XX2407	798.D	<b>DF</b> 1	<b>Analyzed</b> 12/13/18 23:	<b>By</b> 27 CP	<b>Prep D</b> 12/13/1	<b>ate</b> 8 09:00	Prep Batch OP17274	<b>Analytical Batch</b> GXX6552
Run #1 Run #2	<b>Initial</b> 1040 ml		<b>Final Vol</b> 5.0 ml	ume					
PCB List									
CAS No.	Comp	ound		Result	RL	MDL	Units	Q	
12674-11-2 11104-28-2 11141-16-5 53469-21-9 12672-29-6 11097-69-1 11096-82-5	Aroclo Aroclo Aroclo Aroclo Aroclo Aroclo Aroclo	r 1221 r 1232 r 1242 r 1248 r 1254		ND ND ND ND ND ND ND	0.24 0.24 0.24 0.24 0.24 0.24 0.24 0.24	0.094 0.20 0.12 0.11 0.061 0.20 0.073	ug/l ug/l ug/l ug/l ug/l ug/l		
CAS No. 877-09-8	Surrog	gate Rec		<b>Run# 1</b> 19%	Run# 2	Lim	U U		
877-09-8 877-09-8		nloro-m- nloro-m-	•	19% 17%			.66%		

(a) Outside the QC limits. There is no sample left to re-extract.

Decachlorobiphenyl

Decachlorobiphenyl

2051-24-3

2051-24-3

9% a

8% a

J = Indicates an estimated value

10-150%

10-150%

- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound



Client San Lab Samj Matrix: Method: Project:	ple ID: JC7 AQ SW	9316-4 - Ground Wa 846 8260C	ater es, Lodi Street, Syrad	cuse, N	]	Date Sampled: Date Received: Percent Solids:	
Run #1 Run #2	<b>File ID</b> 2V55853.D	<b>DF</b> 1	Analyzed 12/12/18 10:07	<b>By</b> JP	<b>Prep Date</b> n/a	<b>Prep Batc</b> n/a	h Analytical Batch V2V2263
Run #1 Run #2	<b>Purge Volu</b> 5.0 ml	me					

#### **VOA Special List**

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	6.0	ug/l	
71-43-2	Benzene	ND	0.50	0.43	ug/l	
108-86-1	Bromobenzene	ND	1.0	0.55	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	0.82	1.0	0.58	ug/l	J
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>a</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	6.9	ug/l	
104-51-8	n-Butylbenzene	ND	2.0	0.52	ug/l	
135-98-8	sec-Butylbenzene	ND	2.0	0.62	ug/l	
98-06-6	tert-Butylbenzene	ND	2.0	0.69	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.95	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	17.4	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
95-49-8	o-Chlorotoluene	ND	2.0	0.63	ug/l	
106-43-4	p-Chlorotoluene	ND	2.0	0.60	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	1.2	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane <sup>a</sup>	ND	2.0	1.4	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
540-59-0	1,2-Dichloroethene (total)	ND	1.0	0.51	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
142-28-9	1,3-Dichloropropane	ND	1.0	0.43	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound

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JC79316

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Client Sample ID:	MW-5		
Lab Sample ID:	JC79316-4	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	SW846 8260C	Percent Solids:	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

**VOA Special List** 

CAS No.	Compound	Result	RL	MDL	Units	Q
594-20-7	2,2-Dichloropropane	ND	1.0	0.52	ug/l	
563-58-6	1,1-Dichloropropene	ND	1.0	0.82	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
87-68-3	Hexachlorobutadiene	ND	2.0	0.56	ug/l	
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l	
99-87-6	p-Isopropyltoluene	ND	2.0	0.66	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
74-95-3	Methylene bromide	ND	1.0	0.48	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
91-20-3	Naphthalene	ND	5.0	0.98	ug/l	
103-65-1	n-Propylbenzene	ND	2.0	0.60	ug/l	
100-42-5	Styrene	ND	1.0	0.70	ug/l	
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0	0.60	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane <sup>b</sup>	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.90	ug/l	
108-88-3	Toluene	ND	1.0	0.53	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.84	ug/l	
96-18-4	1,2,3-Trichloropropane	ND	2.0	0.70	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	2.0	1.0	ug/l	
108-67-8	1,3,5-Trimethylbenzene	ND	2.0	1.0	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.79	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	ın#1 Run#2 Liı		its	
1868-53-7	Dibromofluoromethane	98%		80-1	20%	
17060-07-0	1,2-Dichloroethane-D4	106%		81-1	24%	
2037-26-5	Toluene-D8	99%		80-1	20%	
460-00-4	4-Bromofluorobenzene	99%	80-120% 80-120%			

(a) Associated CCV outside of control limits low.

(b) This compound in BS is outside in house QC limits bias high.

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

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JC79316

E = Indicates value exceeds calibration range

	<b>Report of Analysis</b>								
Client San Lab Samp Matrix: Method: Project:	le ID: JC7931 AQ - G SW846					Date Sampled: 12/05/18 Date Received: 12/07/18 Percent Solids: n/a			
Run #1 Run #2	<b>File ID</b> XX240799.D	<b>DF</b> 1	<b>Analyzed</b> 12/13/18 23:45	<b>By</b> CP	<b>Prep Date</b> 12/13/18 09:00	Prep Batch OP17274	<b>Analytical Batch</b> GXX6552		
Run #1 Run #2 PCB List	<b>Initial Volume</b> 1040 ml	Final Vo 5.0 ml	lume						

CAS No.	Compound	Result	RL	MDL	Units	Q
12674-11-2 11104-28-2 11141-16-5 53469-21-9	Aroclor 1016 Aroclor 1221 Aroclor 1232 Aroclor 1242	ND ND ND ND	0.24 0.24 0.24 0.24	0.094 0.20 0.12 0.11	ug/l ug/l ug/l ug/l	
12672-29-6 11097-69-1 11096-82-5	Aroclor 1248 Aroclor 1254 Aroclor 1260	ND ND ND	0.24 0.24 0.24	0.061 0.20 0.073	ug/l ug/l ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Lim	its	
877-09-8 877-09-8 2051-24-3 2051-24-3	Tetrachloro-m-xylene Tetrachloro-m-xylene Decachlorobiphenyl Decachlorobiphenyl	133% 121% 63% 61%		11-1 11-1 10-1 10-1	50%	

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound







Client San Lab Samj Matrix: Method: Project:	ple ID: JC A( SV	W-2 79316-5 Q - Ground Wa V846 8260C uanta Resource	ater 28, Lodi Street, Syrad	cuse, N		Date Sampled: Date Received: Percent Solids:	
Run #1 Run #2	<b>File ID</b> 2V55958.I	<b>DF</b> 0 1	<b>Analyzed</b> 12/14/18 11:14	<b>By</b> JTP	<b>Prep Date</b> n/a	<b>Prep Bate</b> n/a	h Analytical Batch V2V2267
Run #1 Run #2	<b>Purge Vol</b> 5.0 ml	ume					

#### **VOA Special List**

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	6.0	ug/l	
71-43-2	Benzene	1.4	0.50	0.43	ug/l	
108-86-1	Bromobenzene	ND	1.0	0.55	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.58	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>a</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	6.9	ug/l	
104-51-8	n-Butylbenzene	ND	2.0	0.52	ug/l	
135-98-8	sec-Butylbenzene	ND	2.0	0.62	ug/l	
98-06-6	tert-Butylbenzene	ND	2.0	0.69	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.95	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	0.97	1.0	0.56	ug/l	J
75-00-3	Chloroethane	0.89	1.0	0.73	ug/l	J
67-66-3	Chloroform	ND	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
95-49-8	o-Chlorotoluene	ND	2.0	0.63	ug/l	
106-43-4	p-Chlorotoluene	ND	2.0	0.60	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	1.2	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	3.1	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	2.5	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	5.9	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	1.4	ug/l	
75-34-3	1,1-Dichloroethane	0.73	1.0	0.57	ug/l	J
107-06-2	1,2-Dichloroethane <sup>b</sup>	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
540-59-0	1,2-Dichloroethene (total)	1.0	1.0	0.51	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
142-28-9	1,3-Dichloropropane	ND	1.0	0.43	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

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JC79316

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Client Sample ID:	MW-2		
Lab Sample ID:	JC79316-5	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	SW846 8260C	<b>Percent Solids:</b>	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

**VOA Special List** 

CAS No.	Compound	Compound Result		MDL	Units	Q
594-20-7	2,2-Dichloropropane	ND	1.0	0.52	ug/l	
563-58-6	1,1-Dichloropropene	ND	1.0	0.82	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
87-68-3	Hexachlorobutadiene	ND	2.0	0.56	ug/l	
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l	
99-87-6	p-Isopropyltoluene	ND	2.0	0.66	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
74-95-3	Methylene bromide	ND	1.0	0.48	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
91-20-3	Naphthalene	ND	5.0	0.98	ug/l	
103-65-1	n-Propylbenzene	ND	2.0	0.60	ug/l	
100-42-5	Styrene	ND	1.0	0.70	ug/l	
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0	0.60	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane <sup>c</sup>	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.90	ug/l	
108-88-3	Toluene	ND	1.0	0.53	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	1.2	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.84	ug/l	
96-18-4	1,2,3-Trichloropropane	ND	2.0	0.70	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	2.0	1.0	ug/l	
108-67-8	1,3,5-Trimethylbenzene	ND	2.0	1.0	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.79	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
CAS No.	Surrogate Recoveries	Run# 1 Run# 2		Lim	its	
1868-53-7	Dibromofluoromethane	99%		80-1	20%	
17060-07-0	1,2-Dichloroethane-D4	109%		81-1	24%	
2037-26-5	Toluene-D8	98%		80-1	20%	
460-00-4	4-Bromofluorobenzene	100%		80-1	20%	

(a) Associated CCV outside of control limits low.

(b) Associated CCV outside of control limits high, sample was ND.

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

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E = Indicates value exceeds calibration range

SGS LabLink@14:59 26-Dec-2018

		Repor	rt of A1	nalysis		Page 3 of 3
Client Sample ID: Lab Sample ID: Matrix: Method: Project:	MW-2 JC79316-5 AQ - Ground Water SW846 8260C Quanta Resources, Lo	odi Street, Sy	yracuse, N	Y	Date Sampled: Date Received: Percent Solids:	
VOA Special List CAS No. Comr	oound	Result	RL	MDL	Units O	

(c) Associated CCV outside of control limits high, sample was ND. This compound in BS is outside in house QC limits bias high.

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



JC79316



2051-24-3

2051-24-3

				Repor	t of An	alysis			Page 1 of 1
Client Sam Lab Sample Matrix: Method: Project:	-	MW-2 JC79316-5 AQ - Ground Water SW846 8082A SW846 3510C Quanta Resources, Lodi Street, Syracuse, NY					Date	e Sampled: Received: ent Solids:	12/05/18 12/07/18 n/a
Run #1 Run #2	File ID XX240		<b>DF</b> 1	<b>Analyzed</b> 12/14/18 01::	<b>By</b> 32 CP	<b>Prep D</b> 12/13/1	<b>ate</b> 8 09:00	Prep Batch OP17274	n <b>Analytical Batch</b> GXX6552
Run #1 Run #2	<b>Initial</b> 1040 m	<b>Volume</b> 1	<b>Final Vol</b> 5.0 ml	ume					
PCB List									
CAS No.	Comp	ound		Result	RL	MDL	Units	Q	
12674-11-2 11104-28-2 11141-16-5 53469-21-9 12672-29-6 11097-69-1 11096-82-5	Aroclo Aroclo Aroclo Aroclo Aroclo	or 1016 or 1221 or 1232 or 1242 or 1248 a or 1254 or 1260		ND ND ND 0.75 1.1 1.2	$\begin{array}{c} 0.24 \\ 0.24 \\ 0.24 \\ 0.24 \\ 0.24 \\ 0.24 \\ 0.24 \\ 0.24 \end{array}$	$\begin{array}{c} 0.094\\ 0.20\\ 0.12\\ 0.11\\ 0.061\\ 0.20\\ 0.073\\ \end{array}$	ug/l ug/l ug/l ug/l ug/l ug/l		
CAS No.	Surro	gate Reco	overies	Run# 1	Run# 2	Lim	its		
877-09-8 877-09-8		hloro-m-2 hloro-m-2	•	141% 113%	11-166% 11-166%				

(a) More than 40 % RPD for detected concentrations between the two GC columns.

41%

46%

Decachlorobiphenyl

Decachlorobiphenyl

J = Indicates an estimated value

10-150%

10-150%

- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





Client San Lab Samj Matrix: Method: Project:	ple ID: JC7 AQ SW8	: MW-1-D JC79316-6 AQ - Ground Water SW846 8260C Quanta Resources, Lodi Street, Syracuse, NY				Date Sampled: 12/05/18 Date Received: 12/07/18 Percent Solids: n/a			
Run #1 Run #2	<b>File ID</b> 2V55873.D	<b>DF</b> 1	<b>Analyzed</b> 12/12/18 17:57	<b>By</b> JP	<b>Prep Date</b> n/a	<b>Prep Batch</b> n/a	Analytical Batch V2V2263		
Run #1 Run #2	<b>Purge Volun</b> 5.0 ml	ne							

#### **VOA Special List**

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	6.0	ug/l	
71-43-2	Benzene	0.52	0.50	0.43	ug/l	
108-86-1	Bromobenzene	ND	1.0	0.55	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.58	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>a</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	6.9	ug/l	
104-51-8	n-Butylbenzene	ND	2.0	0.52	ug/l	
135-98-8	sec-Butylbenzene	ND	2.0	0.62	ug/l	
98-06-6	tert-Butylbenzene	ND	2.0	0.69	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.95	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	0.57	1.0	0.56	ug/l	J
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	ND	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
95-49-8	o-Chlorotoluene	ND	2.0	0.63	ug/l	
106-43-4	p-Chlorotoluene	ND	2.0	0.60	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	1.2	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	0.59	1.0	0.53	ug/l	J
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane <sup>a</sup>	ND	2.0	1.4	ug/l	
75-34-3	1,1-Dichloroethane	2.3	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
540-59-0	1,2-Dichloroethene (total)	1.5	1.0	0.51	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
142-28-9	1,3-Dichloropropane	ND	1.0	0.43	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound



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E = Indicates value exceeds calibration range

Client Sample ID:	MW-1-D		
Lab Sample ID:	JC79316-6	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	SW846 8260C	Percent Solids:	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

**VOA Special List** 

CAS No.	Compound	Result	RL	MDL	Units	Q
594-20-7	2,2-Dichloropropane	ND	1.0	0.52	ug/l	
563-58-6	1,1-Dichloropropene	ND	1.0	0.82	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
87-68-3	Hexachlorobutadiene	ND	2.0	0.56	ug/l	
98-82-8	Isopropylbenzene	ND	1.0	0.65	ug/l	
99-87-6	p-Isopropyltoluene	ND	2.0	0.66	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
74-95-3	Methylene bromide	ND	1.0	0.48	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
91-20-3	Naphthalene	ND	5.0	0.98	ug/l	
103-65-1	n-Propylbenzene	ND	2.0	0.60	ug/l	
100-42-5	Styrene	ND	1.0	0.70	ug/l	
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0	0.60	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane <sup>b</sup>	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.90	ug/l	
108-88-3	Toluene	ND	1.0	0.53	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.84	ug/l	
96-18-4	1,2,3-Trichloropropane	ND	2.0	0.70	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	2.0	1.0	ug/l	
108-67-8	1,3,5-Trimethylbenzene	ND	2.0	1.0	ug/l	
75-01-4	Vinyl chloride	24.0	1.0	0.79	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Lim	its	
1868-53-7	Dibromofluoromethane	95%		80-1	20%	
17060-07-0	1,2-Dichloroethane-D4	104%	81-124%			
2037-26-5	Toluene-D8	99%	80-120%			
460-00-4	4-Bromofluorobenzene	100%		80-1	20%	

(a) Associated CCV outside of control limits low.

(b) This compound in BS is outside in house QC limits bias high.

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound





E = Indicates value exceeds calibration range

53469-21-9 Aroclor 1242

12672-29-6 Aroclor 1248

11097-69-1 Aroclor 1254

11096-82-5 Aroclor 1260

**Surrogate Recoveries** 

Tetrachloro-m-xylene

Tetrachloro-m-xylene

Decachlorobiphenyl

Decachlorobiphenyl

CAS No.

877-09-8

877-09-8

2051-24-3

2051-24-3

			Report	nalysis			Page 1 of 1		
Client Sample ID:MW-1-DLab Sample ID:JC79316-6Matrix:AQ - Ground WaterMethod:SW846 8082ASW846 3510CProject:Quanta Resources, Lodi Street, Syracuse, NY					Date Sampled: 12/05/18 Date Received: 12/07/18 Percent Solids: n/a Y				
Run #1 Run #2	<b>File ID</b> XX240805.D	<b>DF</b> 1	<b>Analyzed</b> 12/14/18 01:50	<b>By</b> CP	<b>Prep D</b> a 12/13/1	ate 8 09:00	Prep Batch OP17274	Analytical Batch GXX6552	
Run #1 Run #2	<b>Initial Volume</b> 1040 ml	Final Vo 5.0 ml	lume						
PCB List									
CAS No.	Compound		Result	RL	MDL	Units	Q		
12674-11-2 11104-28-2 11141-16-5	Aroclor 1221		ND	0.24 0.24 0.24	0.094 0.20 0.12	ug/l ug/l ug/l			

0.24

0.24

0.24

0.24

**Run# 2** 

0.11

0.061

0.20

0.073

Limits

11-166%

11-166%

10-150%

10-150%

ug/l

ug/l

ug/l

ug/l

ND

ND

ND

ND

42%

36%

19%

20%

Run#1

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound



# 4.6

Client Sar Lab Samj Matrix: Method: Project:	ple ID: J( A S	MW-12 JC79316-7 AQ - Ground Water SW846 8260C Quanta Resources, Lodi Street, Syracuse, NY				Date Sampled: 12/05/18 Date Received: 12/07/18 Percent Solids: n/a		
Run #1 Run #2	<b>File ID</b> 2V55872.	<b>DF</b> D 1	<b>Analyzed</b> 12/12/18 17:31	<b>By</b> JP	<b>Prep Date</b> n/a	<b>Prep Batc</b> n/a	h Analytical Batch V2V2263	
Run #1 Run #2	<b>Purge Vo</b> 5.0 ml	lume						

#### **VOA Special List**

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	6.0	ug/l	
71-43-2	Benzene	0.46	0.50	0.43	ug/l	J
108-86-1	Bromobenzene	ND	1.0	0.55	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.58	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>a</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	6.9	ug/l	
104-51-8	n-Butylbenzene	ND	2.0	0.52	ug/l	
135-98-8	sec-Butylbenzene	1.4	2.0	0.62	ug/l	J
98-06-6	tert-Butylbenzene	0.96	2.0	0.69	ug/l	J
75-15-0	Carbon disulfide	ND	2.0	0.95	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	11.2	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	ND	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
95-49-8	o-Chlorotoluene	ND	2.0	0.63	ug/l	
106-43-4	p-Chlorotoluene	ND	2.0	0.60	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	1.2	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	3.4	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	1.8	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane <sup>a</sup>	ND	2.0	1.4	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
540-59-0	1,2-Dichloroethene (total)	ND	1.0	0.51	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
142-28-9	1,3-Dichloropropane	ND	1.0	0.43	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound

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JC79316

E = Indicates value exceeds calibration range

J = Indicates an estimated value

Client Sample ID:	MW-12		
Lab Sample ID:	JC79316-7	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	SW846 8260C	Percent Solids:	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

**VOA Special List** 

CAS No.	Compound	Result	RL	MDL	Units	Q
594-20-7	2,2-Dichloropropane	ND	1.0	0.52	ug/l	
563-58-6	1,1-Dichloropropene	ND	1.0	0.82	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
87-68-3	Hexachlorobutadiene	ND	2.0	0.56	ug/l	
98-82-8	Isopropylbenzene	3.9	1.0	0.65	ug/l	
99-87-6	p-Isopropyltoluene	ND	2.0	0.66	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
74-95-3	Methylene bromide	ND	1.0	0.48	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
91-20-3	Naphthalene	ND	5.0	0.98	ug/l	
103-65-1	n-Propylbenzene	1.3	2.0	0.60	ug/l	J
100-42-5	Styrene	ND	1.0	0.70	ug/l	
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0	0.60	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane <sup>b</sup>	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.90	ug/l	
108-88-3	Toluene	ND	1.0	0.53	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.84	ug/l	
96-18-4	1,2,3-Trichloropropane	ND	2.0	0.70	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	2.0	1.0	ug/l	
108-67-8	1,3,5-Trimethylbenzene	ND	2.0	1.0	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.79	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2 Limits		its	
1868-53-7	Dibromofluoromethane	96%		80-1	20%	
17060-07-0	1,2-Dichloroethane-D4	104%		81-1	24%	
2037-26-5	Toluene-D8	98%	80-120%			
460-00-4	4-Bromofluorobenzene	100%		80-1	20%	

(a) Associated CCV outside of control limits low.

(b) This compound in BS is outside in house QC limits bias high.

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound



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E = Indicates value exceeds calibration range

		Page 1 of 1							
Client Sam Lab Sampl Matrix: Method: Project:	e ID: JC79 AQ SW8	MW-12 JC79316-7 AQ - Ground Water SW846 8270D BY SIM SW846 3510C Quanta Resources, Lodi Street, Syracuse, N				Date Sampled: 12/05/18 Date Received: 12/07/18 Percent Solids: n/a			
Run #1 Run #2	<b>File ID</b> 3P73450.D	<b>DF</b> 1	<b>Analyzed</b> 12/13/18 03:09	By SA	<b>Prep D</b> 12/11/1		<b>Prep Batch</b> OP17214A	<b>Analytical Batch</b> E3P3464	
Run #1 Run #2	<b>Initial Volun</b> 1050 ml	ne Final Vo 1.0 ml	olume						
CAS No.	Compound		Result	RL	MDL	Units	Q		
123-91-1	1,4-Dioxane	2	0.239	0.095	0.046	ug/l			
CAS No.	Surrogate I	Recoveries	<b>Run#</b> 1	Run# 2	Lim	its			
4165-60-0 321-60-8 1718-51-0	Nitrobenzen 2-Fluorobip Terphenyl-d	henyl	73% 62% 45%		29-1 23-1 22-1	22%			

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

E = Indicates value exceeds calibration range

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Client Sar Lab Samp Matrix: Method: Project:	AQ EPA	9316-7 - Ground Wa . 537M BY II	ter D EPA 537 MOD s, Lodi Street, Syrac	cuse, N	Date Sampled:12/05/18Date Received:12/07/18Percent Solids:n/aY			
Run #1 <sup>a</sup> Run #2	<b>File ID</b> 3Q559.D	<b>DF</b> 1	<b>Analyzed</b> 12/18/18 22:49	<b>By</b> AFL	<b>Prep Date</b> 12/14/18 08:45	<b>Prep Batch</b> F:OP73036	<b>Analytical Batch</b> F:S3Q9	
	Initial Volur	ne Final V	olume					

Run #1 250 ml

Run #2

1.0 ml

#### **PFAS List**

CAS No.	Compound	Result	RL	MDL	Units	Q
375-22-4	Perfluorobutanoic acid	4.87	8.0	2.0	ng/l	J
2706-90-3	Perfluoropentanoic acid	ND	4.0	1.5	ng/l	
307-24-4	Perfluorohexanoic acid	1.31	4.0	1.0	ng/l	J
375-85-9	Perfluoroheptanoic acid	ND	2.0	1.0	ng/l	
335-67-1	Perfluorooctanoic acid	3.41	2.0	1.0	ng/l	
375-95-1	Perfluorononanoic acid	ND	2.0	1.0	ng/l	
335-76-2	Perfluorodecanoic acid	ND	4.0	1.0	ng/l	
2058-94-8	Perfluoroundecanoic acid	ND	4.0	1.0	ng/l	
307-55-1	Perfluorododecanoic acid	ND	4.0	1.5	ng/l	
72629-94-8	Perfluorotridecanoic acid	ND	4.0	1.0	ng/l	
376-06-7	Perfluorotetradecanoic acid	ND	4.0	1.0	ng/l	
375-73-5	Perfluorobutanesulfonic acid	ND	2.0	1.0	ng/l	
355-46-4	Perfluorohexanesulfonic acid	ND	2.0	1.0	ng/l	
375-92-8	Perfluoroheptanesulfonic acid	ND	4.0	1.0	ng/l	
1763-23-1	Perfluorooctanesulfonic acid	6.12	2.0	1.5	ng/l	
335-77-3	Perfluorodecanesulfonic acid	ND			ng/l	
754-91-6	PFOSA	ND	4.0	1.0	ng/l	
2355-31-9	MeFOSAA	ND	20	4.0	ng/l	
2991-50-6	EtFOSAA	ND	20	4.0	ng/l	
27619-97-2	6:2 Fluorotelomer sulfonate	ND	8.0	2.0	ng/l	
39108-34-4	8:2 Fluorotelomer sulfonate	ND	8.0	2.0	ng/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Lim	its	
	13C4-PFBA	54%		30-1	40%	
	13C5-PFPeA	59%		40-1	40%	
	13C5-PFHxA	63%		50-1	50%	
	13C4-PFHpA	66%		50-1	50%	
	13C8-PFOA	74%		50-1	50%	
	13C9-PFNA	73%		50-1	50%	
	13C6-PFDA	64%		50-1	50%	
	13C7-PFUnDA	89%		50-1	50%	

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound



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Client Sample ID:	MW-12		
Lab Sample ID:	JC79316-7	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	EPA 537M BY ID EPA 537 MOD	Percent Solids:	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		
-	- ·		

#### **PFAS List**

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
	13C2-PFDoDA 13C2-PFTeDA	78% 75%		50-150% 40-150%
	13C3-PFBS 13C3-PFHxS	58% 59%		50-150% 50-150%
	13C8-PFOS 13C8-FOSA	63% 65%		50-150% 50-150% 30-140%
	d3-MeFOSAA	84%		50-150%
	13C2-6:2FTS 13C2-8:2FTS	91% 78%		50-150% 50-150%

(a) Analysis performed at SGS Orlando, FL.

- J = Indicates an estimated value
- $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$
- N = Indicates presumptive evidence of a compound

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12672-29-6 Aroclor 1248

11097-69-1 Aroclor 1254

11096-82-5 Aroclor 1260

**Surrogate Recoveries** 

Tetrachloro-m-xylene

Tetrachloro-m-xylene

Decachlorobiphenyl

Decachlorobiphenyl

CAS No.

877-09-8

877-09-8

2051-24-3

2051-24-3

			Report	of Ai	nalysis			Page 1 of 1
Client Sam Lab Sampl Matrix: Method: Project:	e ID: JC7931 AQ - G SW846	6-7 round Wate 8082A SV	er W846 3510C Lodi Street, Syrad	cuse, N	īΥ	Date	Received:	12/05/18 12/07/18 n/a
Run #1 Run #2	<b>File ID</b> XX240806.D	<b>DF</b> 1	<b>Analyzed</b> 12/14/18 02:08	<b>By</b> CP	<b>Prep D</b> a 12/13/1	ate 8 09:00	Prep Batch OP17274	<b>Analytical Batch</b> GXX6552
Run #1 Run #2	<b>Initial Volume</b> 1000 ml	<b>Final Vo</b> 5.0 ml	lume					
PCB List								
CAS No.	Compound		Result	RL	MDL	Units	Q	
12674-11-2 11104-28-2 11141-16-5 53469-21-9	Aroclor 1221 Aroclor 1232		ND ND	0.25 0.25 0.25 0.25	0.098 0.21 0.13 0.11	ug/l ug/l ug/l ug/l		

0.25

0.25

0.25

**Run# 2** 

0.063

0.076

Limits

11-166%

11-166%

10-150%

10-150%

0.21

ug/l

ug/l

ug/l

ND

ND

ND

Run#1

119%

98%

50%

54%

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound





Client San Lab Samj Matrix: Method: Project:	ple ID: JC7 AQ SW8	9316-8 - Ground Wa 346 8260C	ater 28, Lodi Street, Syrad	cuse, N	F	2/05/18 2/07/18 ⁄a			
Run #1 Run #2	<b>File ID</b> 2V55871.D	<b>DF</b> 1	<b>Analyzed</b> 12/12/18 17:06	<b>By</b> JP	<b>Prep Date</b> n/a	<b>Prep Batch</b> n/a	<b>Analytical Batch</b> V2V2263		
Run #1 Run #2	<b>Purge Volur</b> 5.0 ml	ne							

**Report of Analysis** 

#### Cull #2

### **VOA Special List**

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	10	6.0	ug/l	
71-43-2	Benzene	2.3	0.50	0.43	ug/l	
108-86-1	Bromobenzene	ND	1.0	0.55	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.48	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.58	ug/l	
75-25-2	Bromoform	ND	1.0	0.63	ug/l	
74-83-9	Bromomethane <sup>a</sup>	ND	2.0	1.6	ug/l	
78-93-3	2-Butanone (MEK)	ND	10	6.9	ug/l	
104-51-8	n-Butylbenzene	0.63	2.0	0.52	ug/l	J
135-98-8	sec-Butylbenzene	1.3	2.0	0.62	ug/l	J
98-06-6	tert-Butylbenzene	ND	2.0	0.69	ug/l	
75-15-0	Carbon disulfide	ND	2.0	0.95	ug/l	
56-23-5	Carbon tetrachloride	ND	1.0	0.55	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.56	ug/l	
75-00-3	Chloroethane	ND	1.0	0.73	ug/l	
67-66-3	Chloroform	ND	1.0	0.50	ug/l	
74-87-3	Chloromethane	ND	1.0	0.76	ug/l	
95-49-8	o-Chlorotoluene	ND	2.0	0.63	ug/l	
106-43-4	p-Chlorotoluene	ND	2.0	0.60	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	2.0	1.2	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.56	ug/l	
106-93-4	1,2-Dibromoethane	ND	1.0	0.48	ug/l	
95-50-1	1,2-Dichlorobenzene	2.6	1.0	0.53	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.54	ug/l	
106-46-7	1,4-Dichlorobenzene	1.1	1.0	0.51	ug/l	
75-71-8	Dichlorodifluoromethane <sup>a</sup>	ND	2.0	1.4	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.57	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.60	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.59	ug/l	
540-59-0	1,2-Dichloroethene (total)	0.93	1.0	0.51	ug/l	J
78-87-5	1,2-Dichloropropane	ND	1.0	0.51	ug/l	
142-28-9	1,3-Dichloropropane	ND	1.0	0.43	ug/l	

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound



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SGS

E = Indicates value exceeds calibration range

# **Report of Analysis**

Client Sample ID:	MW-10		
Lab Sample ID:	JC79316-8	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	SW846 8260C	<b>Percent Solids:</b>	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

**VOA Special List** 

CAS No.	Compound	Result	RL	MDL	Units	Q
594-20-7	2,2-Dichloropropane	ND	1.0	0.52	ug/l	
563-58-6	1,1-Dichloropropene	ND	1.0	0.82	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.47	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.43	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.60	ug/l	
87-68-3	Hexachlorobutadiene	ND	2.0	0.56	ug/l	
98-82-8	Isopropylbenzene	2.3	1.0	0.65	ug/l	
99-87-6	p-Isopropyltoluene	ND	2.0	0.66	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.51	ug/l	
74-95-3	Methylene bromide	ND	1.0	0.48	ug/l	
75-09-2	Methylene chloride	ND	2.0	1.0	ug/l	
91-20-3	Naphthalene	ND	5.0	0.98	ug/l	
103-65-1	n-Propylbenzene	1.9	2.0	0.60	ug/l	J
100-42-5	Styrene	ND	1.0	0.70	ug/l	
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0	0.60	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane <sup>b</sup>	ND	1.0	0.65	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.90	ug/l	
108-88-3	Toluene	ND	1.0	0.53	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	1.0	0.50	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	1.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.54	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.53	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.53	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.84	ug/l	
96-18-4	1,2,3-Trichloropropane	ND	2.0	0.70	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	2.0	1.0	ug/l	
108-67-8	1,3,5-Trimethylbenzene	ND	2.0	1.0	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.79	ug/l	
	m,p-Xylene	ND	1.0	0.78	ug/l	
95-47-6	o-Xylene	ND	1.0	0.59	ug/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Lim	its	
1868-53-7	Dibromofluoromethane	99%		80-1	20%	
17060-07-0	1,2-Dichloroethane-D4	107%		81-1		
2037-26-5	Toluene-D8	99%		80-1		
460-00-4	4-Bromofluorobenzene	100%		80-1		

(a) Associated CCV outside of control limits low.

(b) This compound in BS is outside in house QC limits bias high.

ND = Not detected MDL = Method Detection Limit

RL = Reporting Limit

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

Page 2 of 2

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JC79316

E = Indicates value exceeds calibration range

			Report	of An	alysis			Page 1 of 1
Client Sam Lab Sampl Matrix: Method: Project:	e ID: JC7 AQ SW3	7-10 9316-8 - Ground Wat 846 8270D BY nta Resources,	L	2/05/18 2/07/18 ′a				
Run #1 Run #2	<b>File ID</b> 3P73451.D	<b>DF</b> 1	<b>Analyzed</b> 12/13/18 03:28	By SA	<b>Prep D</b> 12/11/1		<b>Prep Batch</b> OP17214A	<b>Analytical Batch</b> E3P3464
Run #1 Run #2	<b>Initial Volu</b> 1000 ml	ne Final Vo 1.0 ml	lume					
CAS No.	Compound		Result	RL	MDL	Units	Q	
123-91-1	1,4-Dioxan	e	0.225	0.10	0.049	ug/l		
CAS No.	Surrogate	Recoveries	Run# 1	Run# 2	Lim	its		
4165-60-0 321-60-8 1718-51-0	Nitrobenzer 2-Fluorobig Terphenyl-o	henyl	62% 51% 36%		29-1 23-1 22-1	22%		

ND = Not detected MDL = Method Detection Limit RL = Reporting Limit

- E = Indicates value exceeds calibration range
- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound



Run #1 <sup>a</sup>	2Q25416.D	20	12/21/10 21.0/	ALL	12/14/10 00.43	F.UP/3030	F:S2Q394		
	File ID	<b>DF</b> 20	Analyzed 12/21/18 21:07	By AFI	Prep Date 12/14/18 08:45	Prep Batch F:OP73036	e e		
Matrix: Method: Project:	EPA : Quant		ater D EPA 537 MOD s, Lodi Street, Syrac	cuse, N	Perc	Received: ent Solids:			
Lab Samp	nple ID: MW-1 ble ID: JC793				Date	Sampled:	12/05/18		

**Report of Analysis** 

Run #1 260 ml 1.0 ml Run #2

#### **PFAS List**

CAS No.	Compound	Result	RL	M	DL Units	Q
375-22-4	Perfluorobutanoic acid	ND	150	38	ng/l	
2706-90-3	Perfluoropentanoic acid	ND	77	29	ng/l	
307-24-4	Perfluorohexanoic acid	ND	77	19	ng/l	
375-85-9	Perfluoroheptanoic acid	ND	38	19	ng/l	
335-67-1	Perfluorooctanoic acid b	ND	38	19	ng/l	
375-95-1	Perfluorononanoic acid	ND	38	19	ng/l	
335-76-2	Perfluorodecanoic acid	ND	77	19	ng/l	
2058-94-8	Perfluoroundecanoic acid	ND	77	19	ng/l	
307-55-1	Perfluorododecanoic acid	ND	77	29	ng/l	
72629-94-8	Perfluorotridecanoic acid	ND	77	19	ng/l	
376-06-7	Perfluorotetradecanoic acid	ND	77	19	ng/l	
375-73-5	Perfluorobutanesulfonic acid	ND	38	19	ng/l	
355-46-4	Perfluorohexanesulfonic acid	ND	38	19	ng/l	
375-92-8	Perfluoroheptanesulfonic acid	ND	77	19	ng/l	
1763-23-1	Perfluorooctanesulfonic acid	ND	38	29	ng/l	
335-77-3	Perfluorodecanesulfonic acid	ND	77	19	ng/l	
754-91-6	PFOSA	ND	77	19	ng/l	
2355-31-9	MeFOSAA	ND	380	77	ng/l	
2991-50-6	EtFOSAA	ND	380	77	ng/l	
27619-97-2	6:2 Fluorotelomer sulfonate	ND	150	38	ng/l	
39108-34-4	8:2 Fluorotelomer sulfonate	ND	150	38	ng/l	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2		Limits	
	13C4-PFBA	99%			30-140%	
	13C5-PFPeA	101%			40-140%	
	13C5-PFHxA	103%			50-150%	
	13C4-PFHpA	104%			50-150%	
	13C8-PFOA	110%			50-150%	
	13C9-PFNA	110%			50-150%	
	13C6-PFDA	115%			50-150%	
	13C7-PFUnDA	111%			50-150%	

ND = Not detected MDL = Method Detection LimitRL = Reporting LimitE = Indicates value exceeds calibration range  $J = \ Indicates \ an \ estimated \ value$ 

 $B = \ Indicates \ analyte \ found \ in \ associated \ method \ blank$ 

N = Indicates presumptive evidence of a compound



Client Sample ID:	MW-10		
Lab Sample ID:	JC79316-8	Date Sampled:	12/05/18
Matrix:	AQ - Ground Water	Date Received:	12/07/18
Method:	EPA 537M BY ID EPA 537 MOD	Percent Solids:	n/a
Project:	Quanta Resources, Lodi Street, Syracuse, NY		

## **PFAS List**

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
	13C2-PFDoDA	117%		50-150%
	13C2-PFTeDA	124%		40-150%
	13C3-PFBS	97%		50-150%
	13C3-PFHxS	97%		50-150%
	13C8-PFOS	103%		50-150%
	13C8-FOSA	114%		30-140%
	d3-MeFOSAA	107%		50-150%
	13C2-6:2FTS	111%		50-150%
	13C2-8:2FTS	114%		50-150%

(a) Dilution required due to matrix interference. Analysis performed at SGS Orlando, FL.

(b) Associated CCV outside of control limits high, sample was ND.

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound

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Page 2 of 2



			Report	of A	nalysis		Page 1 of 1
Client San Lab Samp Matrix: Method: Project:	le ID: JC7931 AQ - G SW846	6-8 round Wat 8082A S	ter SW846 3510C s, Lodi Street, Syrac	cuse, N	Date Perc	L .	12/05/18 12/07/18 n/a
Run #1 Run #2	<b>File ID</b> XX240807.D	<b>DF</b> 1	<b>Analyzed</b> 12/14/18 02:26	By CP	<b>Prep Date</b> 12/13/18 09:00	Prep Batch OP17274	Analytical Batch GXX6552
Run #1 Run #2	<b>Initial Volume</b> 1040 ml	Final V 5.0 ml	olume				
PCB List							

CAS No.	Compound	Result	RL	MDL	Units	Q
12674-11-2 11104-28-2 11141-16-5 53469-21-9 12672-29-6 11097-69-1 11096-82-5	Aroclor 1016 Aroclor 1221 Aroclor 1232 Aroclor 1242 Aroclor 1248 Aroclor 1254 Aroclor 1260	ND ND ND ND 0.28 ND	0.24 0.24 0.24 0.24 0.24 0.24 0.24 0.24	0.094 0.20 0.12 0.11 0.061 0.20 0.073	ug/l ug/l ug/l ug/l ug/l ug/l ug/l	
<b>CAS No.</b> 877-09-8	Surrogate Recoveries Tetrachloro-m-xylene	<b>Run# 1</b> 95%	Run# 2		its 66%	
877-09-8 2051-24-3 2051-24-3	Tetrachloro-m-xylene Decachlorobiphenyl Decachlorobiphenyl	85% 55% 59%		10-1	66% 50% 50%	

- J = Indicates an estimated value
- B = Indicates analyte found in associated method blank
- N = Indicates presumptive evidence of a compound









Misc. Forms

Custody Documents and Other Forms

Includes the following where applicable:

- Chain of Custody
- Chain of Custody (SGS Orlando, FL)

SGS ACCUTE	6	$,\omega$	CHAI		<b>)F</b> C		OI	ΟY												OF _ ]
ACCUTE	ST		2235 F	toute 130	), Daytor	n, NJ 088					FED-	EX Trackir	g#			Bott	e Order Co	2912	3-15	4316
'			TEL. 732-32		FAX: 7 accutest.c		99/348	50			SGS	Accutest C	iuoto * P <i>l</i>	<u>75_</u> 2	018_4	rq ses	Accutest J	∞# J	TC 7	4316
Client / Reporting Information	Project Name:		Project	Informa	ation							Red	ueste	d Analy	rsis ( se	TEST	CODE	sheet)	1 1	Matrix Codes
Company Name Plum ley Engineering Street Address	(	Juan tu	Res	الات ال	<u>じ</u>								37 Accl)	(00278)						DW - Drinking Water GW - Ground Water WW - Water
8232 Loop Rel		. St				on ( if diffe	rent fro	m Rep	ort to)		_5		37	3						SW - Surface Water SO - Soil
City State Zip Baldwinsville NY 13027 Project Contact E-mail	City S J ra	inse	State	Compar Street A								- are	EPA S							SL- Sludge SED-Sediment OI - Oil LIQ - Other Liquid
Matt Martin	2015	5127									ŝ,		Ū,	Š						AIR - Air SOL - Other Solid
Phone # Fax #	Client Purchase	Order# 5127		City			Sta	əte		Zip	٦d	5 3		2 2 0						WP - Wipe FB-Field Blank
Sampler(s) Name(s) Phone #	Project Manager			Attention	ו:				_		┦ ╹	12	1	6					1	EB-Equipment Blank RB- Rinse Blank
Mutt Murtin Derke Hudson	Dule 1	Wilner	Collection							d Bottles		<u>~</u> اد	1	· .						TB-Trip Blank
SGS Access Service # Field ID / Point of Collection	MEOH/DI Vial #	Date	Time	Sampled by	Matrix	# of bottles	Ŧ	T.	TI	MEOH		PCR	đ	ر ۲						LAB USE ONLY
MW-9		12/5/18	1230	My	GW						X	×	×	X						
1 MW-9 MS			1230	1	1						×	(	X							
mw-9 msi)			1230								X		×							E106
2 Equipment Blunk			1250						$\square$		$\square$		X							503
3 MW-6			13:01					$\square$	$\downarrow$	$\square$	×	×					_			VII32
4 mw-5			13:25	11_				Ш	$\square$	++	<u> </u>	×	_			_	_	<u> </u>		
5 mw-2			13:47	<u>                                     </u>				$\square$	$\square$		×	: x				_	_			1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 - 1997 -
6 mw-1-D			1400		11			$\square$	$\downarrow\downarrow$	++	<u>   ×</u>	( X			_	_				
7 MW-12		-	14:50	11	+			⊥	$\downarrow \downarrow$	++	<u>  ×</u>	×	X	X						
\$ mw-10		×	16:40	V	*			_	$\downarrow$	++	$\left  \right\rangle$	식츠	×	×		-			$\vdash$	
								$\downarrow$	11-	++	$\left  \right $	_			_					
Turnaround Time ( Business days)					040	Data	Deliver	rable In	formati	on			140.00		C	ommen	ts / Spec	al Instru	ictions	
Std. Jø Business Days □ 5 Day RUSH		Accutest PM): / Date			Commer	:ial "A" (L :ial "B" ( L ( Level 3+4	evel 2)		X	NYASP (	ategory A ategory B									needed
3 Day RUSH	INTER	ASESSME	<u>4400</u>	20	NJ Reduc	ed	,			EDD For Other	mat N/S	SAIL	í	1,4	1) . 11	are	(	2,28	'ny/L	- reeded
2 Day RUSH     1 Day RUSH	LABEL	VERIFICATI	ON	口		of Known	-		col Rep	orting				PFA	3			2 ng	12	nexclud
other  Emergency & Rush T/A data available VIA Lablink						Results On sults + QC					C Summa	ry	Sam	ale inv	antory i	verifi		5		Laboratory
Consigning of Suar (A Suar (A Suar )	Sa	mple Custody n	ust be docum	iented	elow eac	belime sa	mples	chang	e poss	estion,	Riciuging	courie	delive	y.	+++					
1 12/6/1	\$ 14:40	Received By:	l	Ľ	1		Relinqui 2	ished B	$\mathbb{Z}$	L	11	7:00		Date Tim	<b>6/1</b> 8	2	aived By:	F	8	
3 Date Time) 3 JZ/7/14	00:00	Beenived By: 3	$\sim$	_			Reling	stied By	. <i>C</i>					Date Tim	e:	Rece 4	eived By:			
Relinquished by: Date Time: 5		Received By:					Custody	y Seal ≇			Intact		Preserv	ed where	applicable			On lee	- 0	cooler Temp.
Form:SM088-01CRev.Date:9/13/16									4											1160-1

Form:SM088-01CRev.Date:9/13/16

JC79316: Chain of Custody Page 1 of 3



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# SGS Sample Receipt Summary

Job Number:	JC79316	Client:		Project:			
Date / Time Received:	12/7/2018 10:00	00 AM Deliver	y Method:	Airbill #'s:			
Cooler Temps (Raw Me Cooler Temps (Co	rrected) °C: Coo		2: (1.8);				
Cooler Security 1. Custody Seals Present: 2. Custody Seals Intact: Cooler Temperature		3. COC Present: 4. Smpl Dates/Time C N_	ики П 1. 2	ample Integrity - Documentation . Sample labels present on bottles: . Container labeling complete: . Sample container label / COC agree:	Y V V	or <u>N</u>	
<ol> <li>Temp criteria achieved</li> <li>Cooler temp verification</li> <li>Cooler media:</li> <li>No. Coolers:</li> </ol>			1.2	ample Integrity - Condition . Sample recvd within HT: . All containers accounted for: . Condition of sample:	V	or N	
Quality Control Preser		<u>N N/A</u> ☑ □	<u>s</u>	ample Integrity - Instructions 1. Analysis requested is clear:		or N	<u>N/A</u>
<ol> <li>2. Trip Blank listed on CO</li> <li>3. Samples preserved pro</li> <li>4. VOCs beadspace free:</li> </ol>	oC:		3	2. Bottles received for unspecified tests 3. Sufficient volume recvd for analysis: 4. Compositing instructions clear:			
•	oc:						V
3. Samples preserved pro	oc:		3 4 5	<ol> <li>Sufficient volume recvd for analysis:</li> <li>Compositing instructions clear:</li> </ol>			

SM089-03 Rev. Date 12/7/17

> JC79316: Chain of Custody Page 2 of 3



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			Job Change Order:	rder:	JC79316
Requested Date:	.: e:	12/11/2018		Received Date:	12/7/2018
Account Name:		Plumley Environmental Engineers	ngineers	Due Date:	12/14/2018
Project Descri	otion:	Project Description: Quanta Resources, Lodi Street, Syracuse, NY	Street, Syracuse, NY	Deliverable:	NYASPB
C/O Initiated By: TF	ed By:	TF PM: TF		TAT (Days):	14
Sample #: JC79316-ALL	IC7931(		Change:		
Dept:		H	Please change to 14 Day TAT. Due 12/21/18.	F. Due 12/21/18.	
TAT:	14				

Date/Time: 12/11/2018 11:43:47 AM To Client: This Change Order is confirmation of the revisions, previously discussed with the Client Service Representative. Above Changes Per: Matthew Martin

JC79316: Chain of Custody Page 3 of 3



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				SGS Nor										CED E	Fracking #					<b>1</b>	der Cort	•	1 of	-		
	JUU			2235 Ro TEL. 732-329																		rol N				
	•					LCom/el		3499	3480					SGS Que	te#					SG8 Jo		J	C7931	16		
	Client / Reporting Information			Project	Inform	ation												Roques	ted Ar	alysis					Matrix Codes	
Compar	iy Name:	Project Name:																							DW - Drinking Water	
			urces, Lodi Str	et, Syracusa,	NY												•								GW - Ground Water	
reet A	odross	Street																							WW - Water SW - Surface Water	
ity	State Zip	CRy		State	Billing I Compan		on (If diffe	rent fi	om Re	port	10)														SO - Soil SI,- Skudge SED-Sediment	
	Contact E-mail nl.Williams@sgs.com	Project#			Street A	ddress																			OI - Oit LIQ - Other Liquid AIR - Air	
hone #		Client Purchase	Order#		City				State		3	İρ				- 1	- 1								SOL - Other Solid WP - Wipe	
																									FB - Field Blank EB-Equipment Blank	
ampler MM	(s) Nemo(s) Pho	ne Project Manager		Collection	Attention	с 1			Manha		reserved E			VY21.											RB - Rinse Blank TB - Trip Blank	
SGS Winple #	Field ID / Point of Collection	MEOH/DI VIII #	Dite	Time	Sampled by	Matrix	# of bottles	ş	HO.	2	40NE Militer	NCOR		LCID537NY21											LAB USE ONLY	
1	MW-9	+	12/5/18	12:30:00 PN		AQ		ŕ		÷+		-1-	H	- X		-	-			-	-	-		$\vdash$		
1D	MW-9 MSD		12/5/18	12:30:00 PN		AQ		+	+	+		╋	+	x									$\left  \right $			
15	MW-9 MS		12/5/18	12:30:00 PN		AQ				+	++	+	╉╍┦	Ŷ				$\rightarrow$				-		-		
2	EQUIPMENT BLANK	-									++	+	++			-	_									
-			12/5/18	12:50:00 PM		AQ		++	-	-	++	+-	+-+	Х		_		_								
7	MW-12		12/5/18	2:50:00 PM	1	AQ		$\square$				-	$\square$	Х	$\rightarrow$	_		_			<u> </u>					
8	MW-10		12/5/18	4:40:00 PM	MM	AQ		$\square$			++		1.1	х												
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	Turnaround Time ( Business days)								erable			-	J						Com	nents /	Specia	Instruc	tions			
	Standard 10 Business Days	Approved By (SGS	PM):/Date:				siai "A" (L siai "8" ( L																		-	
	5 Business Days RUSH						(Level 3+4		9		 			ry e	1											
	3 Business Days RUSH					NJ Reduc	ed			Ē	_ EDC	For	mat .													
	2 Business Days RUSH					Commerc					X Oth	er <u>N</u>	YASF	PΒ	_											
	1 Business Day EMERGENCY X Other 7						Commerc					mar												,		
Eme	rgency & Rush T/A data svallable via Aplink Appn	wal needed for RUS	WEmergency TAT				Commerc	ial °C'	= Res	its +	QC Sum	mary	+ Part	iaí Raw e	inte					htt	p://ww	w.sgs.	com/er	derms	-and-conditions	
Rella		(Time: 😙	Sample Cus Received By:	ody must be d			v each tin		mples quished								y. Me/Tim		_	Receive	1	<u> </u>	-7	Ļ		-
		in giti		Fec	( E	<u></u>		2	qual NC		1	Fe	e d		Eχ	ľ	eve / 110	e,		2 4	C/	-	$/ \zeta$	-	12/12/1	3'
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Reline	ulshed by:		Received By: 5					Custo	ody Seal	#				ntect kot intect	Pr At	eserved	Aftere a	plicable	herm (D			on too	,	Cooler 1	femp. 'C	

JC79316.xis Rev. Dste: 4/10/18

> JC79316: Chain of Custody Page 1 of 3 SGS Orlando, FL



SGS

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# SGS Sample Receipt Summary

Job Number: JC7931	6	Client:	ALNJ		Project:	QUANTA RE	SOU	RCES		
Date / Time Received: 12/12/20	018 9:30:00 A	M	Delivery Method:	FED EX	Airbill #'s	: <u>100189177</u>	71660	00328	311004	461413714308
Therm ID: IR 1;			Therm CF: -0.2;			# of Coolers	<b>s:</b> 1			
Cooler Temps (Raw Measure	d) °C: Coole	er 1: (4.0	D);							
Cooler Temps (Correcte	d) °C: Coole	er 1: (3.8	3);							
Cooler Information	Y or	N		Sample Information			Y	or	N	N/A
1. Custody Seals Present	$\checkmark$			1. Sample labels present	on bottles		$\checkmark$			
2. Custody Seals Intact	$\checkmark$			2. Samples preserved pro	operly		$\checkmark$			
3. Temp criteria achieved	$\checkmark$			3. Sufficient volume/conta	ainers recvd	for analysis:	$\checkmark$			
4. Cooler temp verification	IR Gun			4. Condition of sample			Intac	<u>zt</u>		
5. Cooler media	Ice (Bag)			5. Sample recvd within H	т		$\checkmark$			
				6. Dates/Times/IDs on C	OC match Sa	imple Label	$\checkmark$			
Trip Blank Information	Y or	<u>N</u>	N/A	7. VOCs have headspace	e					
1. Trip Blank present / cooler				8. Bottles received for un	specified tes	ts			$\checkmark$	
2. Trip Blank listed on COC				9. Compositing instruction	ns clear					$\checkmark$
	W or	e	N/A	10. Voa Soil Kits/Jars rec	eived past 4	Bhrs?				$\checkmark$
				11. % Solids Jar received	1?					$\checkmark$
3. Type Of TB Received				12. Residual Chlorine Pre	esent?					$\checkmark$
Misc. Information										
Number of Encores: 25-Gran	n	5-Gram	Num	ber of 5035 Field Kits:		Number of La	b Filte	red Me	etals:	
Test Strip Lot #s:	pH 0-3	23031	5pH	H 10-12 219813A	_	Other: (Speci	ify)			
Residual Chlorine Test Strip Lo	t #:									
Comments										
Comments										

SM001 Rev. Date 05/24/17

Technician: SHAYLAP Date: 12/12/2018 9:30:00 A

 Reviewer:
 SP
 Date:
 12/12/2018

JC79316: Chain of Custody Page 2 of 3



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		Job Change	e Order:	JC79316
Requested Date:	12/11/2018		Received Date:	12/7/2018
Account Name:	Plumley Env	vironmental Engineers	Due Date:	12/14/2018
Project Description	n: Quanta Res	ources, Lodi Street, Syracuse, NY	Deliverable:	NYASPB
C/O Initiated	By: TF	PM: TF	TAT (Days):	14
Sample #: JC7 Dept: TAT: 14	9316-ALL	Change: Please change to 14 Day	TAT. Due 12/21/18.	

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JC79316: Chain of Custody

Above Changes Per: Matthew Martin

Date/Time: 12/11/2018 11:45:51 AM

Page 3 of 3

To Client: This Change Order is confirmation of the revisions, previously discussed with the Client Service Representative.

Page 1 of 1



