

The order of this data package is as follows:

1. Chain-of-Custody/Lab Request
2. Copies of COCs
3. Validation Report
4. Laboratory analysis

Comments:

[illegible]

[illegible]

LANL SMO Los Alamos NM		Chain of Custody/Analysis Request															COC/Lab Request #: 2018-1803 Page 1 of 1						
Client Contact:		Lab Agreement #:			Site Name: Los Alamos National Laboratory															Rad Screening Info: Lab Reporting Limit Type: Method Detection Limit			
		Project Number:			MSGP-Hg	WSP-8260B-VOA	WSP-8270C-SVOA	WSP-8330B-NMED HEXMOD	WSP-All Metals	WSP-CN(T)	WSP-GENINORG+PerChlorate	WSP-GrossA/B	WSP-LL-H-3	WSP-NH3+NO3/NO2+PO4	WSP-RAD	WSP-TKN+TOC							
		Analysis Turnaround Time:																					
		24 Hour - <input type="checkbox"/> Other - <input checked="" type="checkbox"/>																					
		7 Days - <input type="checkbox"/>																					
		14 Days - <input type="checkbox"/>																					
		21 Days - <input type="checkbox"/>																					
		28 Days - <input type="checkbox"/>																					
Field Sample ID	Sample Date	Sample Time	Sample Matrix																				
CAWA-18-55	Feb 21 2018	10:44	W					1		1				1									
CAWA-18-112	Feb 21 2018	10:44	W		21																		
CAWA-18-56	Feb 21 2018	10:44	W	1	2	2	3		1		1	1		1	1								
Special Instructions:																							
Relinquished by: Tanya VanderVis				Print Name: Tanya VanderVis				Date/Time: 2-21-18 1340				Received by: Renee Onstott				Print Name: Renee Onstott				Date/Time: 2/21/18 1340			
Relinquished by:				Print Name:				Date/Time:				Received by:				Print Name:				Date/Time:			
Relinquished by:				Print Name:				Date/Time:				Received by:				Print Name:				Date/Time:			

[illegible]

Sampling Plan ID/Name: 11669 (R-26 P2-2)COC: 2018-1811

TEST – Explosives		YES	NO
Samples collected from a WFO area? (TAs -08, 09, 11, 14, 15, 16, 22, 36, 37, 39, 40, and 49)			X
Field Test for Explosives Results		YES	NO
HE SPOT test result positive. If YES - Do not transport.			X

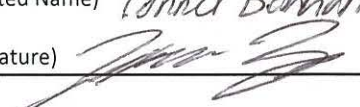
TEST – Chemical Preservation		YES	NO
Samples are chemically preserved?		X	X
Field Team Member Statement		YES	NO
Chemical preservation exceeds limits given 40 CFR 136, Table II – Required Containers, Preservation Techniques and Holding Times (footnote 3). If YES - Do not ship.			X

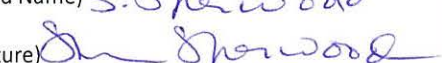
TEST – Field Screen			YES	NO
The sample has field screening measurements of alpha activity and beta activity?				X
Sample Activity (dpm/100cm ²)	Shipment Activity (dpm*g/100cm ²)	Sampled Location	YES	NO
Alpha detectable	AND Alpha ≥ 160,000	AT TA-1 and adjacent hillsides, TA-21, Acid Canyon, MDA C at TA-50, Area G at TA-54, TA-48, or TA-49		X
Alpha ≥ 125	AND Alpha ≥ 1,250,000	AT other locations		
Beta ≥ 1,500	AND Beta ≥ 15,000,000	AT any location		
The sample Alpha ≥ 16,000,000 dpm*g/100cm ² or Beta ≥ 160,000,000 dpm*g/100cm ² . If YES – Do not ship.				
On the external surface of the sample container, alpha activity ≥ 24 dpm/cm ² , beta activity ≥ 240 dpm/cm ² , or surface activity ≥ 0.5 mR/hr. If YES – Do not ship.				
The sample is tentatively identified as DOT Hazard Class 7 (Radioactive). The shipment is labeled Radioactive Material, Excepted Package – Limited Quantity of Material – UN2910, based on field screening measurements of alpha and beta activity.				

TEST - Location			YES	NO
Prior analytical measurements of radioactive isotopes are available?			X	
Sample Activity (pCi/g)	Shipment Activity (pCi)		YES	NO
• Am-241 ≥ 27 pCi/g	AND	Am-241 ≥ 270,000 pCi Total		X
• Cs-137 ≥ 270 pCi/g	AND	Cs-137 ≥ 270,000 pCi Total		
• Pu-238 ≥ 27 pCi/g	AND	Pu-238 ≥ 270,000 pCi Total		
• Pu-239/240 ≥ 27 pCi/g	AND	Pu-239/240 ≥ 270,000 pCi Total		
• Th-228 ≥ 27 pCi/g	AND	Th-228 ≥ 270,000 pCi Total		
• U-234 ≥ 270 pCi/g	AND	U-234 ≥ 1,600,000,000 pCi Total		
• U-238 ≥ 270 pCi/g	AND	U-238 ≥ unlimited		
• H-3 ≥ 27,000,000 pCi/g	AND	H-3 ≥ 27,000,000,000 pCi Total		
Am-241, Pu-238, Pu-239/240, or Th-228 ≥ 27,000,000 pCi; or Cs-137 ≥ 270,000,000 pCi or U-234 ≥ 160,000,000 pCi; or H-3 ≥ 1 Ci. If YES – Do not ship.				
The sample is tentatively identified as DOT Hazard Class 7 (Radioactive). The shipment is labeled Radioactive Material, Excepted Package – Limited Quantity of Material – UN2910, based on prior analytical measurements of radioactive isotopes.				

TEST – AK		YES	NO
The shippers documented knowledge of the sample positively identifies appropriate labeling.			X
Documented Field Team Member Statement		YES	NO
The sample is tentatively identified as DOT Hazard Class 7 (Radioactive). The shipment is labeled <i>Radioactive Material, Excepted Package – Limited Quantity of Material – UN2910</i> , and the sample is submitted to ARS or RP for hazard classification analysis.			X

These samples do not meet the criteria for classification in any hazard class according to regulation OSHA 29 CFR 1910.1200. The sample(s) contained in this shipment have been assigned a tentative proper DOT shipping name, hazard class, identification number, and packing group, based on the shipper's knowledge of the sample:

Hazard Assessment Completed By:	Date/Time
(Printed Name) <u>Tanner Barham</u>	<u>2-21-2018</u>
(Signature) 	<u>1300</u>

Hazard Assessment Reviewed By:	Date/Time
(Printed Name) <u>S. Sherwood</u>	<u>2/21/18</u>
(Signature) 	<u>13:00</u>

Sampling Plan ID/Name: 11669 Bulldog Spring, R-26 SICOC: 2018-1811

TEST – Explosives		YES	NO
Samples collected from a WFO area? (TAs -08, 09, 11, 14, 15, 16, 22, 36, 37, 39, 40, and 49)		<input checked="" type="checkbox"/>	<input type="checkbox"/>
Field Test for Explosives Results		YES	NO
HE SPOT test result positive. If YES - Do not transport.		<input checked="" type="checkbox"/>	<input type="checkbox"/>

TEST – Chemical Preservation		YES	NO
Samples are chemically preserved?		<input checked="" type="checkbox"/>	<input type="checkbox"/>
Field Team Member Statement		YES	NO
Chemical preservation exceeds limits given 40 CFR 136, Table II – Required Containers, Preservation Techniques and Holding Times (footnote 3). If YES - Do not ship.		<input checked="" type="checkbox"/>	<input type="checkbox"/>

TEST – Field Screen			YES	NO
The sample has field screening measurements of alpha activity and beta activity?			<input checked="" type="checkbox"/>	<input type="checkbox"/>
Sample Activity (dpm/100cm ²)	Shipment Activity (dpm*g/100cm ²)	Sampled Location	YES	NO
Alpha detectable	AND Alpha ≥ 160,000	AT TA-1 and adjacent hillsides, TA-21, Acid Canyon, MDA C at TA-50, Area G at TA-54, TA-48, or TA-49	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Alpha ≥ 125	AND Alpha ≥ 1,250,000	AT other locations	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Beta ≥ 1,500	AND Beta ≥ 15,000,000	AT any location	<input checked="" type="checkbox"/>	<input type="checkbox"/>
The sample Alpha ≥ 16,000,000 dpm*g/100cm ² or Beta ≥ 160,000,000 dpm*g/100cm ² . If YES – Do not ship.			<input checked="" type="checkbox"/>	<input type="checkbox"/>
On the external surface of the sample container, alpha activity ≥ 24 dpm/cm ² , beta activity ≥ 240 dpm/cm ² , or surface activity ≥ 0.5 mR/hr. If YES – Do not ship.			<input checked="" type="checkbox"/>	<input type="checkbox"/>
The sample is tentatively identified as DOT Hazard Class 7 (Radioactive). The shipment is labeled Radioactive Material, Excepted Package – Limited Quantity of Material – UN2910, based on field screening measurements of alpha and beta activity.			<input checked="" type="checkbox"/>	<input type="checkbox"/>

TEST - Location			YES	NO
Prior analytical measurements of radioactive isotopes are available?			<input checked="" type="checkbox"/>	<input type="checkbox"/>
Sample Activity (pCi/g)	Shipment Activity (pCi)		YES	NO
• Am-241 ≥ 27 pCi/g	AND	Am-241 ≥ 270,000 pCi Total	<input checked="" type="checkbox"/>	<input type="checkbox"/>
• Cs-137 ≥ 270 pCi/g	AND	Cs-137 ≥ 270,000 pCi Total	<input checked="" type="checkbox"/>	<input type="checkbox"/>
• Pu-238 ≥ 27 pCi/g	AND	Pu-238 ≥ 270,000 pCi Total	<input checked="" type="checkbox"/>	<input type="checkbox"/>
• Pu-239/240 ≥ 27 pCi/g	AND	Pu-239/240 ≥ 270,000 pCi Total	<input checked="" type="checkbox"/>	<input type="checkbox"/>
• Th-228 ≥ 27 pCi/g	AND	Th-228 ≥ 270,000 pCi Total	<input checked="" type="checkbox"/>	<input type="checkbox"/>
• U-234 ≥ 270 pCi/g	AND	U-234 ≥ 1,600,000,000 pCi Total	<input checked="" type="checkbox"/>	<input type="checkbox"/>
• U-238 ≥ 270 pCi/g	AND	U-238 ≥ unlimited	<input checked="" type="checkbox"/>	<input type="checkbox"/>
• H-3 ≥ 27,000,000 pCi/g	AND	H-3 ≥ 27,000,000,000 pCi Total	<input checked="" type="checkbox"/>	<input type="checkbox"/>
Am-241, Pu-238, Pu-239/240, or Th-228 ≥ 27,000,000 pCi; or Cs-137 ≥ 270,000,000 pCi or U-234 ≥ 160,000,000 pCi; or H-3 ≥ 1 Ci. If YES – Do not ship.			<input checked="" type="checkbox"/>	<input type="checkbox"/>
The sample is tentatively identified as DOT Hazard Class 7 (Radioactive). The shipment is labeled Radioactive Material, Excepted Package – Limited Quantity of Material – UN2910, based on prior analytical measurements of radioactive isotopes.			<input checked="" type="checkbox"/>	<input type="checkbox"/>

TEST – AK		YES	NO
The shippers documented knowledge of the sample positively identifies appropriate labeling.		<input checked="" type="checkbox"/>	<input type="checkbox"/>
Documented Field Team Member Statement		YES	NO
The sample is tentatively identified as DOT Hazard Class 7 (Radioactive). The shipment is labeled <i>Radioactive Material, Excepted Package – Limited Quantity of Material – UN2910</i> , and the sample is submitted to ARS or RP for hazard classification analysis.		<input checked="" type="checkbox"/>	<input type="checkbox"/>

These samples do not meet the criteria for classification in any hazard class according to regulation OSHA 29 CFR 1910.1200. The sample(s) contained in this shipment have been assigned a tentative proper DOT shipping name, hazard class, identification number, and packing group, based on the shipper's knowledge of the sample:

Hazard Assessment Completed By:	Date/Time
(Printed Name) <u>Tanya VanderVis</u>	<u>2-21-18</u>
(Signature) <u>Tanya VanderVis</u>	<u>1340</u>

Hazard Assessment Reviewed By:	Date/Time
(Printed Name) <u>Ranee Onstott</u>	<u>2/21/18</u>
(Signature) <u>Ranee Onstott</u>	<u>1340</u>

DATA VALIDATION REPORT

Chain Of Custody No. 2018-1811

1. Distribution Of Samples In EDD.

SDG	Analytical Method	Regular Samples	Field Duplicates	Trip Blanks	Field Blanks	Equipment Blanks
444566	EPA:120.1	2	1			
444566	EPA:150.1	2	1			
444566	EPA:160.1	2	1			
444566	EPA:170.0	5	2	3		
444566	EPA:245.2	4	2			
444566	EPA:300.0	2	1			
444566	EPA:310.1	2	1			
444566	EPA:335.4	2	1			
444566	EPA:350.1	2	1			
444566	EPA:351.2	2	1			
444566	EPA:353.2	2	1			
444566	EPA:365.4	2	1			
444566	EPA:900	2	1			
444566	EPA:901.1	2	1			
444566	EPA:905.0	2	1			
444566	HASL-300:AM-241	2	1			
444566	HASL-300:ISOPU	2	1			
444566	HASL-300:ISOU	2	1			
444566	SM:A2340B	2	1			
444566	SW-846:6010C	2	1			
444566	SW-846:6020	2	1			
444566	SW-846:6850	2	1			
444566	SW-846:8260B	3	1	3		
444566	SW-846:8270D	3	1			
444566	SW-846:8330B	3	1			
444566	SW-846:9060	2	1			

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DATA VALIDATION REPORT

SDG	Analytical Method	Analysis Lot ID	Prep Lot ID	Regular Samples	Field Duplicates	Trip Blanks	Field Blanks	Equipment Blanks	Method Blanks	Matrix Spikes	Matrix Spike Dups	Analytical Spikes	Post-Digestion Spikes	Lab Control Samples	Lab Control Sample Dups	Blank Spike	Blank Spike Dups	Lab Duplicates	Storage Blanks	Preparation Blanks	Reagent Blanks
444566	EPA:120.1	1742552	1742552	2	1									1			1				
444566	EPA:150.1	1741049	1741049	2	1									1			1				
444566	EPA:160.1	1742467	1742467	2	1				1					1			1				
444566	EPA:170.0	NA	NA	5	2	3															
444566	EPA:245.2	1742549	1742548	4	2				1	1				1			1				
444566	EPA:300.0	1741949	1741949	2	1				1					1			1				
444566	EPA:310.1	1741047	1741047	2	1					2				1			2				
444566	EPA:335.4	1740518	1740517	2	1				1	1				1			1				
444566	EPA:350.1	1741276	1741272	2	1				1	1				1			1				
444566	EPA:351.2	1742074	1742073	2	1				1	2				1			2				
444566	EPA:353.2	1742109	1742109	2	1				1					1			1				
444566	EPA:365.4	1742097	1742096	2	1				1	1				1			1				
444566	EPA:900	1742280	1742280	2	1				1	1	1			1			1				
444566	EPA:901.1	1741844	1741844	2	1				1					1			1				
444566	EPA:905.0	1742296	1742296	2	1				1	1				1			1				
444566	HASL-300:AM-241	1741983	1741983	2	1				1					1			1				
444566	HASL-300:ISOPU	1741984	1741984	2	1				1					1			1				
444566	HASL-300:ISOU	1741985	1741985	2	1				1					1			1				
444566	SM:A2340B	1747180	1747180	2	1																
444566	SW-846:6010C	1741857	1741856	2	1				1	1				1			1				
444566	SW-846:6020	1741864	1741863	2	1				1	1				1			1				
444566	SW-846:6850	1742510	1742509	2	1				1	1	1			1							
444566	SW-846:8260B	1744716	1744716	3	1	3			2					4							
444566	SW-846:8270D	1742249	1742247	3	1				1	1	1			1							
444566	SW-846:8330B	1742521	1742520	3	1				1	1	1			1							
444566	SW-846:9060	1741854	1741854	2	1				1					1			1				

2. Distribution Of Analytes In EDD.

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DATA VALIDATION REPORT

Analytical Method	Analytical Method Category	Field Sample ID	Lab Sample ID	Sample Purpose	Target Analytes	Surrogates	Spiked Compounds	TICS
EPA:120.1	GENERAL CHEMISTRY	CAMO-18-16	1203979623	DUP	1	0	0	0
EPA:120.1	GENERAL CHEMISTRY	CAPA-18-1	444566001	REG	1	0	0	0
EPA:120.1	GENERAL CHEMISTRY	CAPA-18-10	444566005	FD	1	0	0	0
EPA:120.1	GENERAL CHEMISTRY	CAWA-18-55	444566011	REG	1	0	0	0
EPA:120.1	GENERAL CHEMISTRY	LCS	1203979622	LCS	0	0	1	0
EPA:150.1	GENERAL CHEMISTRY	CAMO-18-1	1203976495	DUP	1	0	0	0
EPA:150.1	GENERAL CHEMISTRY	CAPA-18-1	444566001	REG	1	0	0	0
EPA:150.1	GENERAL CHEMISTRY	CAPA-18-10	444566005	FD	1	0	0	0
EPA:150.1	GENERAL CHEMISTRY	CAWA-18-55	444566011	REG	1	0	0	0
EPA:150.1	GENERAL CHEMISTRY	LCS	1203976494	LCS	0	0	1	0
EPA:160.1	GENERAL CHEMISTRY	CAMO-18-19	1203979423	DUP	1	0	0	0
EPA:160.1	GENERAL CHEMISTRY	CAPA-18-1	444566001	REG	1	0	0	0
EPA:160.1	GENERAL CHEMISTRY	CAPA-18-10	444566005	FD	1	0	0	0
EPA:160.1	GENERAL CHEMISTRY	CAWA-18-55	444566011	REG	1	0	0	0
EPA:160.1	GENERAL CHEMISTRY	LCS	1203979422	LCS	0	0	1	0
EPA:160.1	GENERAL CHEMISTRY	MB	1203979421	MB	1	0	0	0
EPA:170.0	VOC	CAPA-18-1	444566001	REG	1	0	0	0
EPA:170.0	VOC	CAPA-18-10	444566005	FD	1	0	0	0
EPA:170.0	VOC	CAPA-18-12	444566006	FD	1	0	0	0
EPA:170.0	VOC	CAPA-18-2	444566002	REG	1	0	0	0
EPA:170.0	VOC	CAPA-18-6	444566004	FTB	1	0	0	0
EPA:170.0	VOC	CAWA-18-111	444566010	FTB	1	0	0	0
EPA:170.0	VOC	CAWA-18-112	444566014	FTB	1	0	0	0
EPA:170.0	VOC	CAWA-18-54	444566008	REG	1	0	0	0
EPA:170.0	VOC	CAWA-18-55	444566011	REG	1	0	0	0
EPA:170.0	VOC	CAWA-18-56	444566012	REG	1	0	0	0
EPA:245.2	INORGANIC	CAPA-18-1	1203979606	DUP	1	0	0	0
EPA:245.2	INORGANIC	CAPA-18-1	1203979607	MS	0	0	1	0
EPA:245.2	INORGANIC	CAPA-18-1	444566001	REG	1	0	0	0
EPA:245.2	INORGANIC	CAPA-18-10	444566005	FD	1	0	0	0
EPA:245.2	INORGANIC	CAPA-18-12	444566006	FD	1	0	0	0
EPA:245.2	INORGANIC	CAPA-18-2	444566002	REG	1	0	0	0
EPA:245.2	INORGANIC	CAWA-18-55	444566011	REG	1	0	0	0
EPA:245.2	INORGANIC	CAWA-18-56	444566012	REG	1	0	0	0
EPA:245.2	INORGANIC	LCS	1203979605	LCS	0	0	1	0
EPA:245.2	INORGANIC	MB	1203979604	MB	1	0	0	0
EPA:300.0	GENERAL CHEMISTRY	CAPA-18-1	444566001	REG	4	0	0	0
EPA:300.0	GENERAL CHEMISTRY	CAPA-18-10	444566005	FD	4	0	0	0
EPA:300.0	GENERAL CHEMISTRY	CAWA-18-55	444566011	REG	4	0	0	0

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DATA VALIDATION REPORT

Analytical Method	Analytical Method Category	Field Sample ID	Lab Sample ID	Sample Purpose	Target Analytes	Surrogates	Spiked Compounds	TICS
EPA:300.0	GENERAL CHEMISTRY	CTUA-17-151327	1203978238	DUP	4	0	0	0
EPA:300.0	GENERAL CHEMISTRY	LCS	1203978237	LCS	0	0	4	0
EPA:300.0	GENERAL CHEMISTRY	MB	1203978236	MB	4	0	0	0
EPA:310.1	GENERAL CHEMISTRY	CAMO-18-1	1203976486	DUP	2	0	0	0
EPA:310.1	GENERAL CHEMISTRY	CAMO-18-1	1203976488	MS	0	0	1	0
EPA:310.1	GENERAL CHEMISTRY	CAPA-18-1	444566001	REG	2	0	0	0
EPA:310.1	GENERAL CHEMISTRY	CAPA-18-10	444566005	FD	2	0	0	0
EPA:310.1	GENERAL CHEMISTRY	CAWA-18-29	1203976487	DUP	2	0	0	0
EPA:310.1	GENERAL CHEMISTRY	CAWA-18-29	1203976489	MS	0	0	1	0
EPA:310.1	GENERAL CHEMISTRY	CAWA-18-55	444566011	REG	2	0	0	0
EPA:310.1	GENERAL CHEMISTRY	LCS	1203976485	LCS	0	0	1	0
EPA:335.4	INORGANIC	CAPA-18-12	444566006	FD	1	0	0	0
EPA:335.4	INORGANIC	CAPA-18-2	444566002	REG	1	0	0	0
EPA:335.4	INORGANIC	CAWA-18-28	1203975197	DUP	1	0	0	0
EPA:335.4	INORGANIC	CAWA-18-28	1203975198	MS	0	0	1	0
EPA:335.4	INORGANIC	CAWA-18-56	444566012	REG	1	0	0	0
EPA:335.4	INORGANIC	LCS	1203975196	LCS	0	0	1	0
EPA:335.4	INORGANIC	MB	1203975195	MB	1	0	0	0
EPA:350.1	GENERAL CHEMISTRY	CAPA-18-1	1203976872	DUP	1	0	0	0
EPA:350.1	GENERAL CHEMISTRY	CAPA-18-1	1203976874	MS	0	0	1	0
EPA:350.1	GENERAL CHEMISTRY	CAPA-18-1	444566001	REG	1	0	0	0
EPA:350.1	GENERAL CHEMISTRY	CAPA-18-10	444566005	FD	1	0	0	0
EPA:350.1	GENERAL CHEMISTRY	CAWA-18-55	444566011	REG	1	0	0	0
EPA:350.1	GENERAL CHEMISTRY	LCS	1203976870	LCS	0	0	1	0
EPA:350.1	GENERAL CHEMISTRY	MB	1203976869	MB	1	0	0	0
EPA:351.2	GENERAL CHEMISTRY	CAPA-18-12	444566006	FD	1	0	0	0
EPA:351.2	GENERAL CHEMISTRY	CAPA-18-2	444566002	REG	1	0	0	0
EPA:351.2	GENERAL CHEMISTRY	CAWA-18-28	1203978466	DUP	1	0	0	0
EPA:351.2	GENERAL CHEMISTRY	CAWA-18-28	1203978467	MS	0	0	1	0
EPA:351.2	GENERAL CHEMISTRY	CAWA-18-36	1203978468	DUP	1	0	0	0
EPA:351.2	GENERAL CHEMISTRY	CAWA-18-36	1203978469	MS	0	0	1	0
EPA:351.2	GENERAL CHEMISTRY	CAWA-18-56	444566012	REG	1	0	0	0
EPA:351.2	GENERAL CHEMISTRY	LCS	1203978465	LCS	0	0	1	0
EPA:351.2	GENERAL CHEMISTRY	MB	1203978464	MB	1	0	0	0
EPA:353.2	GENERAL CHEMISTRY	CAPA-18-1	444566001	REG	1	0	0	0
EPA:353.2	GENERAL CHEMISTRY	CAPA-18-10	444566005	FD	1	0	0	0
EPA:353.2	GENERAL CHEMISTRY	CAWA-18-27	1203978569	DUP	1	0	0	0
EPA:353.2	GENERAL CHEMISTRY	CAWA-18-55	444566011	REG	1	0	0	0
EPA:353.2	GENERAL CHEMISTRY	LCS	1203978567	LCS	0	0	1	0

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DATA VALIDATION REPORT

Analytical Method	Analytical Method Category	Field Sample ID	Lab Sample ID	Sample Purpose	Target Analytes	Surrogates	Spiked Compounds	TICS
EPA:353.2	GENERAL CHEMISTRY	MB	1203978566	MB	1	0	0	0
EPA:365.4	GENERAL CHEMISTRY	CAPA-18-1	444566001	REG	1	0	0	0
EPA:365.4	GENERAL CHEMISTRY	CAPA-18-10	444566005	FD	1	0	0	0
EPA:365.4	GENERAL CHEMISTRY	CAWA-18-27	1203978531	DUP	1	0	0	0
EPA:365.4	GENERAL CHEMISTRY	CAWA-18-27	1203978532	MS	0	0	1	0
EPA:365.4	GENERAL CHEMISTRY	CAWA-18-55	444566011	REG	1	0	0	0
EPA:365.4	GENERAL CHEMISTRY	LCS	1203978530	LCS	0	0	1	0
EPA:365.4	GENERAL CHEMISTRY	MB	1203978529	MB	1	0	0	0
EPA:900	RAD	CAPA-18-12	444566006	FD	2	0	0	0
EPA:900	RAD	CAPA-18-2	444566002	REG	2	0	0	0
EPA:900	RAD	CAWA-18-40	1203978997	DUP	2	0	0	0
EPA:900	RAD	CAWA-18-40	1203978998	MS	0	0	2	0
EPA:900	RAD	CAWA-18-40	1203978999	MSD	0	0	2	0
EPA:900	RAD	CAWA-18-56	444566012	REG	2	0	0	0
EPA:900	RAD	LCS	1203979000	LCS	0	0	2	0
EPA:900	RAD	MB	1203978996	MB	2	0	0	0
EPA:901.1	RAD	CAPA-18-12	444566006	FD	5	0	0	0
EPA:901.1	RAD	CAPA-18-2	1203978022	DUP	5	0	0	0
EPA:901.1	RAD	CAPA-18-2	444566002	REG	5	0	0	0
EPA:901.1	RAD	CAWA-18-56	444566012	REG	5	0	0	0
EPA:901.1	RAD	LCS	1203978023	LCS	0	0	3	0
EPA:901.1	RAD	MB	1203978021	MB	5	0	0	0
EPA:905.0	RAD	CAPA-18-12	444566006	FD	1	0	0	0
EPA:905.0	RAD	CAPA-18-2	444566002	REG	1	0	0	0
EPA:905.0	RAD	CAWA-18-56	444566012	REG	1	0	0	0
EPA:905.0	RAD	CAWA-18-67	1203979042	DUP	1	0	0	0
EPA:905.0	RAD	CAWA-18-67	1203979043	MS	0	0	1	0
EPA:905.0	RAD	LCS	1203979044	LCS	0	0	1	0
EPA:905.0	RAD	MB	1203979041	MB	1	0	0	0
HASL-300:AM-241	RAD	CAPA-18-12	444566006	FD	1	0	0	0
HASL-300:AM-241	RAD	CAPA-18-2	1203978299	DUP	1	0	0	0
HASL-300:AM-241	RAD	CAPA-18-2	444566002	REG	1	0	0	0
HASL-300:AM-241	RAD	CAWA-18-56	444566012	REG	1	0	0	0
HASL-300:AM-241	RAD	LCS	1203978300	LCS	0	0	1	0
HASL-300:AM-241	RAD	MB	1203978298	MB	1	0	0	0
HASL-300:ISOPU	RAD	CAPA-18-12	444566006	FD	2	0	0	0
HASL-300:ISOPU	RAD	CAPA-18-2	1203978302	DUP	2	0	0	0
HASL-300:ISOPU	RAD	CAPA-18-2	444566002	REG	2	0	0	0
HASL-300:ISOPU	RAD	CAWA-18-56	444566012	REG	2	0	0	0

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Analytical Method	Analytical Method Category	Field Sample ID	Lab Sample ID	Sample Purpose	Target Analytes	Surrogates	Spiked Compounds	TICS
HASL-300:ISOPU	RAD	LCS	1203978303	LCS	0	0	1	0
HASL-300:ISOPU	RAD	MB	1203978301	MB	2	0	0	0
HASL-300:ISOU	RAD	CAPA-18-12	444566006	FD	3	0	0	0
HASL-300:ISOU	RAD	CAPA-18-2	1203978306	DUP	3	0	0	0
HASL-300:ISOU	RAD	CAPA-18-2	444566002	REG	3	0	0	0
HASL-300:ISOU	RAD	CAWA-18-56	444566012	REG	3	0	0	0
HASL-300:ISOU	RAD	LCS	1203978307	LCS	0	0	1	0
HASL-300:ISOU	RAD	MB	1203978305	MB	3	0	0	0
SM:A2340B	INORGANIC	CAPA-18-1	444566001	REG	1	0	0	0
SM:A2340B	INORGANIC	CAPA-18-10	444566005	FD	1	0	0	0
SM:A2340B	INORGANIC	CAWA-18-55	444566011	REG	1	0	0	0
SW-846:6010C	INORGANIC	CAPA-18-1	1203978046	DUP	17	0	0	0
SW-846:6010C	INORGANIC	CAPA-18-1	1203978047	MS	0	0	17	0
SW-846:6010C	INORGANIC	CAPA-18-1	444566001	REG	17	0	0	0
SW-846:6010C	INORGANIC	CAPA-18-10	444566005	FD	17	0	0	0
SW-846:6010C	INORGANIC	CAWA-18-55	444566011	REG	17	0	0	0
SW-846:6010C	INORGANIC	LCS	1203978045	LCS	0	0	17	0
SW-846:6010C	INORGANIC	MB	1203978044	MB	17	0	0	0
SW-846:6020	INORGANIC	CAPA-18-1	1203978056	DUP	11	0	0	0
SW-846:6020	INORGANIC	CAPA-18-1	1203978057	MS	0	0	11	0
SW-846:6020	INORGANIC	CAPA-18-1	444566001	REG	11	0	0	0
SW-846:6020	INORGANIC	CAPA-18-10	444566005	FD	11	0	0	0
SW-846:6020	INORGANIC	CAWA-18-55	444566011	REG	11	0	0	0
SW-846:6020	INORGANIC	LCS	1203978055	LCS	0	0	11	0
SW-846:6020	INORGANIC	MB	1203978054	MB	11	0	0	0
SW-846:6850	LCMS/MS PERCHLORATE	CAPA-18-1	444566001	REG	1	0	0	0
SW-846:6850	LCMS/MS PERCHLORATE	CAPA-18-10	444566005	FD	1	0	0	0
SW-846:6850	LCMS/MS PERCHLORATE	CAWA-18-27	1203979481	MS	0	0	1	0
SW-846:6850	LCMS/MS PERCHLORATE	CAWA-18-27	1203979482	MSD	0	0	1	0
SW-846:6850	LCMS/MS PERCHLORATE	CAWA-18-55	444566011	REG	1	0	0	0
SW-846:6850	LCMS/MS PERCHLORATE	LCS	1203979480	LCS	0	0	1	0
SW-846:6850	LCMS/MS PERCHLORATE	MB	1203979479	MB	1	0	0	0
SW-846:8260B	VOC	CAPA-18-12	444566006	FD	80	3	0	0
SW-846:8260B	VOC	CAPA-18-2	444566002	REG	80	3	0	0
SW-846:8260B	VOC	CAPA-18-6	444566004	FTB	80	3	0	0
SW-846:8260B	VOC	CAWA-18-111	444566010	FTB	80	3	0	0
SW-846:8260B	VOC	CAWA-18-112	444566014	FTB	80	3	0	0
SW-846:8260B	VOC	CAWA-18-54	444566009	REG	80	3	0	0
SW-846:8260B	VOC	CAWA-18-56	444566012	REG	80	3	0	0

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Analytical Method	Analytical Method Category	Field Sample ID	Lab Sample ID	Sample Purpose	Target Analytes	Surrogates	Spiked Compounds	TICS
SW-846:8260B	VOC	LCS	1203984433	LCS	0	3	70	0
SW-846:8260B	VOC	LCS	1203984434	LCS	0	3	10	0
SW-846:8260B	VOC	LCS	1203985581	LCS	0	3	70	0
SW-846:8260B	VOC	LCS	1203985582	LCS	0	3	10	0
SW-846:8260B	VOC	MB	1203984431	MB	80	3	0	0
SW-846:8260B	VOC	MB	1203985580	MB	80	3	0	0
SW-846:8270D	SVOC	CAPA-18-12	444566006	FD	80	6	0	0
SW-846:8270D	SVOC	CAPA-18-2	444566002	REG	80	6	0	0
SW-846:8270D	SVOC	CAWA-18-54	444566009	REG	80	6	0	0
SW-846:8270D	SVOC	CAWA-18-56	1203978894	MS	0	6	76	0
SW-846:8270D	SVOC	CAWA-18-56	1203978895	MSD	0	6	76	0
SW-846:8270D	SVOC	CAWA-18-56	444566012	REG	80	6	0	0
SW-846:8270D	SVOC	LCS	1203978893	LCS	0	6	76	0
SW-846:8270D	SVOC	MB	1203978892	MB	80	6	0	0
SW-846:8330B	LCMS/MS HIGH	CAPA-18-12	444566007	FD	23	1	0	0
SW-846:8330B	LCMS/MS HIGH	CAPA-18-2	1203979522	MS	0	1	23	0
SW-846:8330B	LCMS/MS HIGH	CAPA-18-2	1203979523	MSD	0	1	23	0
SW-846:8330B	LCMS/MS HIGH	CAPA-18-2	444566003	REG	23	1	0	0
SW-846:8330B	LCMS/MS HIGH	CAWA-18-54	444566008	REG	23	1	0	0
SW-846:8330B	LCMS/MS HIGH	CAWA-18-56	444566013	REG	23	1	0	0
SW-846:8330B	LCMS/MS HIGH	LCS	1203979521	LCS	0	1	23	0
SW-846:8330B	LCMS/MS HIGH	MB	1203979520	MB	23	1	0	0
SW-846:9060	GENERAL CHEMISTRY	CAMO-18-28	1203980265	DUP	1	0	0	0
SW-846:9060	GENERAL CHEMISTRY	CAPA-18-12	444566006	FD	1	0	0	0
SW-846:9060	GENERAL CHEMISTRY	CAPA-18-2	444566002	REG	1	0	0	0
SW-846:9060	GENERAL CHEMISTRY	CAWA-18-56	444566012	REG	1	0	0	0
SW-846:9060	GENERAL CHEMISTRY	LCS	1203980263	LCS	0	0	1	0
SW-846:9060	GENERAL CHEMISTRY	MB	1203980262	MB	1	0	0	0

3. Are any analytes missing?

No.

4. Were any holding times exceeded?

No.

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5. Any contaminants in blanks?

						Blank Lab Result	Lab Qualifier	Blank Lab Units	Blank Lab Detection Limit
Blank FS ID	Blank Lab Sample	Blank Type	Analytical Method	Sample	Parameter Name				
CAPA-18-6	444566004	TRIP BLANK	EPA:170.0	W	Temperature	2		Deg C	
CAWA-18-111	444566010	TRIP BLANK	EPA:170.0	W	Temperature	2		Deg C	
CAWA-18-112	444566014	TRIP BLANK	EPA:170.0	W	Temperature	2		Deg C	

No.

6. Any surrogate recoveries outside the control limits?

Field Sample ID	Lab Sample ID	Analytical Method	Parameter Name	Analysis Lot ID	Analysis Date	Spike Recovery	Upper Limit	Lower Limit	Rejection Limit
CAWA-18-54	444566009	SW-846:8270D	4-Terphenyl-d14	1742249	02-27-2018	25	121	36	10
CAWA-18-54	444566009	SW-846:8270D	Fluorobiphenyl[2-]	1742249	02-27-2018	21	112	32	10
CAWA-18-54	444566009	SW-846:8270D	Nitrobenzene-d5	1742249	02-27-2018	24	115	36	10
CAWA-18-54	444566009	SW-846:8270D	Phenol-d5	1742249	02-27-2018	12	91	15	10
CAWA-18-54	444566009	SW-846:8270D	Tribromophenol[2,4,6-]	1742249	02-27-2018	24	124	32	10

7. Any MS/MSD recoveries or RPDs outside the control limits?

No.

8. Any LCS/LCSD or BS/BSD recoveries or RPDs outside the control limits?

DATA VALIDATION REPORT

LCS Lab Sample	LCSD Lab	Analytical Method	Parameter Name	Lab Lot ID	Analysis	Sample Matrix	LCS Spike Recovery	LCSD Spike Recovery	Upper Limit	Lower Limit	Upper Rejection Limit	Lower Rejection Limit	RPD	RPD Limit
1203985581		SW-846:8260B	Hexachlorobutadiene	1744716	03-07-2018	W	66		136	72		10		

9. Any Field Duplicate RPDs outside the desired limits?

No.

10. Any Lab Duplicate RPDs outside the desired limits?

No.

11. Any required reporting limits exceeded?

No.

12. Additional Validator's Comments.

13. Display Flagged Data.

Location ID	COC Number	Field Sample ID	Sample Purpose	Analysis Type Code	Analytical Suite	Analytical Method	Parameter Name	Lab Qualifier	Validation Qualifier	Validation Reason Codes	Detect Flag	Lab Result	Lab Units	Report Result	Report Units	Report MDA	Report Uncertainty	Lab Matrix	Sample Date	Percent	Analysis Lot ID	Validation Status Code	Use Flag
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	HASL-300:AM-241	Americium-241	U	U	R5	N	0.00371	pCi/L	0.00371	pCi/L	0.0313	0.00586	W	02/21/2018		1741983	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	EPA:901.1	Cesium-137	U	U	R5	N	-0.267	pCi/L	-0.267	pCi/L	3.35	0.935	W	02/21/2018		1741844	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	EPA:901.1	Cobalt-60	U	U	R5	N	0.428	pCi/L	0.428	pCi/L	3.36	0.794	W	02/21/2018		1741844	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	EPA:900	Gross alpha	U	U	R5	N	1.07	pCi/L	1.07	pCi/L	2.66	0.781	W	02/21/2018		1742280	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	VOC	SW-846:8260B	Hexachlorobutadiene	U	U	V12a	N	0.300	ug/L	0.300	ug/L			W	02/21/2018		1744716	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	EPA:901.1	Neptunium-237	U	U	R5	N	-0.991	pCi/L	-0.991	pCi/L	6.54	1.82	W	02/21/2018		1741844	VAL	Y

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Location ID	COC Number	Field Sample ID	Sample Purpose	Analysis Type Code	Analytical Suite	Analytical Method	Parameter Name	Lab Qualifier	Validation Qualifier	Validation Reason Codes	Detect Flag	Lab Result	Lab Units	Report Result	Report Units	Report MDA	Report Uncertainty	Lab Matrix	Sample Date	Percent	Analysis Lot ID	Validation Status Code	Use Flag
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	HASL-300:ISOPU	Plutonium-238	U	U	R5	N	0.00425	pCi/L	0.00425	pCi/L	0.0369	0.00672	W	02/21/2018		1741984	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	HASL-300:ISOPU	Plutonium-239/240	U	U	R5	N	0.00213	pCi/L	0.00213	pCi/L	0.0297	0.00766	W	02/21/2018		1741984	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	EPA:901.1	Potassium-40	U	U	R5	N	-19.6	pCi/L	-19.6	pCi/L	47.8	13.5	W	02/21/2018		1741844	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	EPA:901.1	Sodium-22	U	U	R5	N	-0.544	pCi/L	-0.544	pCi/L	2.54	0.680	W	02/21/2018		1741844	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	EPA:905.0	Strontium-90	U	U	R5	N	0.016	pCi/L	0.016	pCi/L	0.278	0.0812	W	02/21/2018		1742296	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-12	FD	INIT	RAD	HASL-300:ISOU	Uranium-235/236	U	U	R5	N	0.0153	pCi/L	0.0153	pCi/L	0.0641	0.00811	W	02/21/2018		1741985	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	HASL-300:AM-241	Americium-241	U	U	R5	N	0.0071	pCi/L	0.0071	pCi/L	0.030	0.00561	W	02/21/2018		1741983	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	EPA:901.1	Cesium-137	U	U	R5	N	-0.858	pCi/L	-0.858	pCi/L	3.61	1.14	W	02/21/2018		1741844	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	EPA:901.1	Cobalt-60	U	U	R5	N	0.0705	pCi/L	0.0705	pCi/L	4.29	1.11	W	02/21/2018		1741844	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	EPA:900	Gross alpha	U	U	R5	N	0.947	pCi/L	0.947	pCi/L	2.78	0.785	W	02/21/2018		1742280	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	VOC	SW-846:8260B	Hexachlorobutadiene	U	UJ	V12a	N	0.300	ug/L	0.300	ug/L			W	02/21/2018		1744716	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	EPA:901.1	Neptunium-237	U	U	R5	N	-0.856	pCi/L	-0.856	pCi/L	6.58	1.85	W	02/21/2018		1741844	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	HASL-300:ISOPU	Plutonium-238	U	U	R5	N	0.00332	pCi/L	0.00332	pCi/L	0.0577	0.00742	W	02/21/2018		1741984	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	HASL-300:ISOPU	Plutonium-239/240	U	U	R5	N	0.0166	pCi/L	0.0166	pCi/L	0.0464	0.011	W	02/21/2018		1741984	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	EPA:901.1	Potassium-40	U	U	R5	N	-19.9	pCi/L	-19.9	pCi/L	58.0	16.7	W	02/21/2018		1741844	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	EPA:901.1	Sodium-22	U	U	R5	N	-2.1	pCi/L	-2.1	pCi/L	3.40	1.13	W	02/21/2018		1741844	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	EPA:905.0	Strontium-90	U	U	R5	N	-0.168	pCi/L	-0.168	pCi/L	0.482	0.131	W	02/21/2018		1742296	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-2	REG	INIT	RAD	HASL-300:ISOU	Uranium-235/236	U	U	R5	N	0.0209	pCi/L	0.0209	pCi/L	0.0625	0.0123	W	02/21/2018		1741985	VAL	Y
Bulldog Spring	2018-1811	CAPA-18-6	FTB	INIT	VOC	SW-846:8260B	Hexachlorobutadiene	U	UJ	V12a	N	0.300	ug/L	0.300	ug/L			W	02/21/2018		1744716	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-111	FTB	INIT	VOC	SW-846:8260B	Hexachlorobutadiene	U	UJ	V12a	N	0.300	ug/L	0.300	ug/L			W	02/21/2018		1744716	VAL	Y
R-26 S1	2018-1811	CAWA-18-112	FTB	INIT	VOC	SW-846:8260B	Hexachlorobutadiene	U	UJ	V12a	N	0.300	ug/L	0.300	ug/L			W	02/21/2018		1744716	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Acenaphthene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Acenaphthylene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Aniline	U	UJ	SV3a	N	4.33	ug/L	4.33	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Anthracene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Atrazine	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Azobenzene	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Benzidine	U	UJ	SV3a	N	4.02	ug/L	4.02	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Benzo(a)anthracene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Benzo(a)pyrene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Benzo(b)fluoranthene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018		1742249	VAL	Y

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Location ID	COC Number	Field Sample ID	Sample Purpose	Analysis Type Code	Analytical Suite	Analytical Method	Paramter Name	Lab Qualifier	Validation Qualifier	Validation Reason Codes	Detect Flag	Lab Result	Lab Units	Report Result	Report Units	Report MDA	Report Uncertainty	Lab Matrix	Sample Date	Percent	Analysis Lot ID	Validation Status Code	Use Flag
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Benzo(g,h,i)perylene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Benzo(k)fluoranthene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Benzoic Acid	U	UJ	SV3a	N	6.19	ug/L	6.19	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Benzyl Alcohol	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Bis(2-chloroethoxy)methane	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Bis(2-chloroethyl)ether	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Bis(2-ethylhexyl)phthalate	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Bromophenyl-phenylether[4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Butylbenzylphthalate	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Chloro-3-methylphenol[4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Chloroaniline[4-]	U	UJ	SV3a	N	3.40	ug/L	3.40	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Chloronaphthalene[2-]	U	UJ	SV3a	N	0.423	ug/L	0.423	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Chlorophenol[2-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Chlorophenyl-phenyl[4-] Ether	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Chrysene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Di-n-butylphthalate	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Di-n-octylphthalate	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dibenz(a,h)anthracene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dibenzofuran	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dichlorobenzene[1,2-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dichlorobenzene[1,3-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dichlorobenzene[1,4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dichlorobenzidine[3,3'-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dichlorophenol[2,4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Diethylphthalate	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dimethyl Phthalate	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dimethylphenol[2,4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dinitro-2-methylphenol[4,6-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dinitrophenol[2,4-]	U	UJ	SV3a	N	5.15	ug/L	5.15	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dinitrotoluene[2,4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dinitrotoluene[2,6-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	

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R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dinoseb	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Dioxane[1,4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Diphenylamine	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Fluoranthene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Fluorene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Hexachlorobenzene	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	VOC	SW-846:8260B	Hexachlorobutadiene	U	UJ	V12a	N	0.300	ug/L	0.300	ug/L			W	02/21/2018	1744716	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Hexachlorobutadiene	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Hexachlorocyclopentadiene	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Hexachloroethane	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Indeno(1,2,3-cd)pyrene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Isophorone	U	UJ	SV3a	N	3.61	ug/L	3.61	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Methylnaphthalene[1-]	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Methylnaphthalene[2-]	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Methylphenol[2-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Methylphenol[3-,4-]	U	UJ	SV3a	N	3.81	ug/L	3.81	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Naphthalene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitroaniline[2-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitroaniline[3-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitroaniline[4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitrobenzene	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitrophenol[2-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitrophenol[4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitroso-di-n-butylamine[N-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitroso-di-n-propylamine[N-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitrosodiethylamine[N-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitrosodimethylamine[N-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Nitrosopyrrolidine[N-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Oxybis(1-chloropropane)[2,2'-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Pentachlorobenzene	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Pentachlorophenol	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018	1742249	VAL	Y	

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R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Phenanthrene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Phenol	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Pyrene	U	UJ	SV3a	N	0.309	ug/L	0.309	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Pyridine	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Tetrachlorobenzene[1,2,4,5]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Tetrachlorophenol[2,3,4,6-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Trichlorobenzene[1,2,4-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Trichlorophenol[2,4,5-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 PZ-2	2018-1811	CAWA-18-54	REG	INIT	SVOC	SW-846:8270D	Trichlorophenol[2,4,6-]	U	UJ	SV3a	N	3.09	ug/L	3.09	ug/L			W	02/21/2018		1742249	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	HASL-300:AM-241	Americium-241	U	U	R5	N	0.00213	pCi/L	0.00213	pCi/L	0.036	0.00369	W	02/21/2018		1741983	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	EPA:901.1	Cesium-137	U	U	R5	N	0.577	pCi/L	0.577	pCi/L	3.51	0.918	W	02/21/2018		1741844	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	EPA:901.1	Cobalt-60	U	U	R5	N	-1.13	pCi/L	-1.13	pCi/L	3.37	0.961	W	02/21/2018		1741844	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	EPA:900	Gross alpha	U	U	R5	N	0.7	pCi/L	0.7	pCi/L	2.76	0.765	W	02/21/2018		1742280	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	VOC	SW-846:8260B	Hexachlorobutadiene	U	UJ	V12a	N	0.300	ug/L	0.300	ug/L			W	02/21/2018		1744716	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	EPA:901.1	Neptunium-237	U	U	R5	N	0.767	pCi/L	0.767	pCi/L	7.56	2.40	W	02/21/2018		1741844	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	HASL-300:ISOPU	Plutonium-238	U	U	R5	N	-0.0065	pCi/L	-0.0065	pCi/L	0.0377	0.00782	W	02/21/2018		1741984	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	HASL-300:ISOPU	Plutonium-239/240	U	U	R5	N	0.0065	pCi/L	0.0065	pCi/L	0.0303	0.0108	W	02/21/2018		1741984	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	EPA:901.1	Potassium-40	U	U	R5	N	27.8	pCi/L	27.8	pCi/L	38.3	20.4	W	02/21/2018		1741844	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	EPA:901.1	Sodium-22	U	U	R5	N	0.267	pCi/L	0.267	pCi/L	3.82	0.951	W	02/21/2018		1741844	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	EPA:905.0	Strontium-90	U	U	R5	N	-0.0293	pCi/L	-0.0293	pCi/L	0.241	0.0694	W	02/21/2018		1742296	VAL	Y
R-26 S1	2018-1811	CAWA-18-56	REG	INIT	RAD	HASL-300:ISOU	Uranium-235/236	U	U	R5	N	0.0244	pCi/L	0.0244	pCi/L	0.0853	0.0129	W	02/21/2018		1741985	VAL	Y

Reason Code

Description

J_LAB

The analytical laboratory qualified the detected result as estimated (J) because the result was less the PQL but greater than the MDL

NQ

The analytical laboratory did not qualify the analyte as not detected and/or any other standard qualifire. The analyte is detected in the sample.

R5

Analyte is not detected because the amount reported is less than the MDC.

SV3a

The surrogate is < the Lower Acceptance Level (LAL) but >=10%R. Follow the external laboratory limits located within the associated data package.

U_LAB

The analytical laboratory qualified the analyte as not detected.

Only results shown in Section 13 'Display Flagged Data' are current as of this report generation. All other sections are valid for the date the COC data was inserted into EIM, and may have changed due to data updates in the intervening time.

DATA VALIDATION REPORT

Reason Code

Description

V12a

The LCS percent recovery was < the LAL but >10%. Follow the external laboratory limits located within the associated data package.

14. Usable Result Count.

Field Sample ID	Location ID	Sample Purpose	Analytical Method	No. Unuseable Records	Total Records
CAPA-18-1	Bulldog Spring	REG	EPA:120.1	0	1
CAPA-18-1	Bulldog Spring	REG	EPA:150.1	0	1
CAPA-18-1	Bulldog Spring	REG	EPA:160.1	0	1
CAPA-18-1	Bulldog Spring	REG	EPA:170.0	0	1
CAPA-18-1	Bulldog Spring	REG	EPA:245.2	0	1
CAPA-18-1	Bulldog Spring	REG	EPA:300.0	0	4
CAPA-18-1	Bulldog Spring	REG	EPA:310.1	0	2
CAPA-18-1	Bulldog Spring	REG	EPA:350.1	0	1
CAPA-18-1	Bulldog Spring	REG	EPA:353.2	0	1
CAPA-18-1	Bulldog Spring	REG	EPA:365.4	0	1
CAPA-18-1	Bulldog Spring	REG	SM:A2340B	0	1
CAPA-18-1	Bulldog Spring	REG	SW-846:6010C	0	17
CAPA-18-1	Bulldog Spring	REG	SW-846:6020	0	11
CAPA-18-1	Bulldog Spring	REG	SW-846:6850	0	1
CAPA-18-10	Bulldog Spring	FD	EPA:120.1	0	1
CAPA-18-10	Bulldog Spring	FD	EPA:150.1	0	1
CAPA-18-10	Bulldog Spring	FD	EPA:160.1	0	1
CAPA-18-10	Bulldog Spring	FD	EPA:170.0	0	1
CAPA-18-10	Bulldog Spring	FD	EPA:245.2	0	1
CAPA-18-10	Bulldog Spring	FD	EPA:300.0	0	4
CAPA-18-10	Bulldog Spring	FD	EPA:310.1	0	2
CAPA-18-10	Bulldog Spring	FD	EPA:350.1	0	1
CAPA-18-10	Bulldog Spring	FD	EPA:353.2	0	1
CAPA-18-10	Bulldog Spring	FD	EPA:365.4	0	1
CAPA-18-10	Bulldog Spring	FD	SM:A2340B	0	1
CAPA-18-10	Bulldog Spring	FD	SW-846:6010C	0	17
CAPA-18-10	Bulldog Spring	FD	SW-846:6020	0	11
CAPA-18-10	Bulldog Spring	FD	SW-846:6850	0	1
CAPA-18-12	Bulldog Spring	FD	EPA:170.0	0	1
CAPA-18-12	Bulldog Spring	FD	EPA:245.2	0	1
CAPA-18-12	Bulldog Spring	FD	EPA:335.4	0	1

Only results shown in Section 13 'Display Flagged Data' are current as of this report generation. All other sections are valid for the date the COC data was inserted into EIM, and may have changed due to data updates in the intervening time.

DATA VALIDATION REPORT

Field Sample ID	Location ID	Sample Purpose	Analytical Method	No. Unuseable Records	Total Records
CAPA-18-12	Bulldog Spring	FD	EPA:351.2	0	1
CAPA-18-12	Bulldog Spring	FD	EPA:900	0	2
CAPA-18-12	Bulldog Spring	FD	EPA:901.1	0	5
CAPA-18-12	Bulldog Spring	FD	EPA:905.0	0	1
CAPA-18-12	Bulldog Spring	FD	HASL-300:AM-241	0	1
CAPA-18-12	Bulldog Spring	FD	HASL-300:ISOPU	0	2
CAPA-18-12	Bulldog Spring	FD	HASL-300:ISOU	0	3
CAPA-18-12	Bulldog Spring	FD	SW-846:8260B	0	80
CAPA-18-12	Bulldog Spring	FD	SW-846:8270D	0	80
CAPA-18-12	Bulldog Spring	FD	SW-846:8330B	0	23
CAPA-18-12	Bulldog Spring	FD	SW-846:9060	0	1
CAPA-18-2	Bulldog Spring	REG	EPA:170.0	0	1
CAPA-18-2	Bulldog Spring	REG	EPA:245.2	0	1
CAPA-18-2	Bulldog Spring	REG	EPA:335.4	0	1
CAPA-18-2	Bulldog Spring	REG	EPA:351.2	0	1
CAPA-18-2	Bulldog Spring	REG	EPA:900	0	2
CAPA-18-2	Bulldog Spring	REG	EPA:901.1	0	5
CAPA-18-2	Bulldog Spring	REG	EPA:905.0	0	1
CAPA-18-2	Bulldog Spring	REG	HASL-300:AM-241	0	1
CAPA-18-2	Bulldog Spring	REG	HASL-300:ISOPU	0	2
CAPA-18-2	Bulldog Spring	REG	HASL-300:ISOU	0	3
CAPA-18-2	Bulldog Spring	REG	SW-846:8260B	0	80
CAPA-18-2	Bulldog Spring	REG	SW-846:8270D	0	80
CAPA-18-2	Bulldog Spring	REG	SW-846:8330B	0	23
CAPA-18-2	Bulldog Spring	REG	SW-846:9060	0	1
CAPA-18-6	Bulldog Spring	FTB	EPA:170.0	0	1
CAPA-18-6	Bulldog Spring	FTB	SW-846:8260B	0	80
CAWA-18-111	R-26 PZ-2	FTB	EPA:170.0	0	1
CAWA-18-111	R-26 PZ-2	FTB	SW-846:8260B	0	80
CAWA-18-112	R-26 S1	FTB	EPA:170.0	0	1
CAWA-18-112	R-26 S1	FTB	SW-846:8260B	0	80
CAWA-18-54	R-26 PZ-2	REG	EPA:170.0	0	1
CAWA-18-54	R-26 PZ-2	REG	SW-846:8260B	0	80
CAWA-18-54	R-26 PZ-2	REG	SW-846:8270D	0	80
CAWA-18-54	R-26 PZ-2	REG	SW-846:8330B	0	23
CAWA-18-55	R-26 S1	REG	EPA:120.1	0	1

Only results shown in Section 13 'Display Flagged Data' are current as of this report generation. All other sections are valid for the date the COC data was inserted into EIM, and may have changed due to data updates in the intervening time.

DATA VALIDATION REPORT

Field Sample ID	Location ID	Sample Purpose	Analytical Method	No. Unuseable Records	Total Records
CAWA-18-55	R-26 S1	REG	EPA:150.1	0	1
CAWA-18-55	R-26 S1	REG	EPA:160.1	0	1
CAWA-18-55	R-26 S1	REG	EPA:170.0	0	1
CAWA-18-55	R-26 S1	REG	EPA:245.2	0	1
CAWA-18-55	R-26 S1	REG	EPA:300.0	0	4
CAWA-18-55	R-26 S1	REG	EPA:310.1	0	2
CAWA-18-55	R-26 S1	REG	EPA:350.1	0	1
CAWA-18-55	R-26 S1	REG	EPA:353.2	0	1
CAWA-18-55	R-26 S1	REG	EPA:365.4	0	1
CAWA-18-55	R-26 S1	REG	SM:A2340B	0	1
CAWA-18-55	R-26 S1	REG	SW-846:6010C	0	17
CAWA-18-55	R-26 S1	REG	SW-846:6020	0	11
CAWA-18-55	R-26 S1	REG	SW-846:6850	0	1
CAWA-18-56	R-26 S1	REG	EPA:170.0	0	1
CAWA-18-56	R-26 S1	REG	EPA:245.2	0	1
CAWA-18-56	R-26 S1	REG	EPA:335.4	0	1
CAWA-18-56	R-26 S1	REG	EPA:351.2	0	1
CAWA-18-56	R-26 S1	REG	EPA:900	0	2
CAWA-18-56	R-26 S1	REG	EPA:901.1	0	5
CAWA-18-56	R-26 S1	REG	EPA:905.0	0	1
CAWA-18-56	R-26 S1	REG	HASL-300:AM-241	0	1
CAWA-18-56	R-26 S1	REG	HASL-300:ISOPU	0	2
CAWA-18-56	R-26 S1	REG	HASL-300:ISOU	0	3
CAWA-18-56	R-26 S1	REG	SW-846:8260B	0	80
CAWA-18-56	R-26 S1	REG	SW-846:8270D	0	80
CAWA-18-56	R-26 S1	REG	SW-846:8330B	0	23
CAWA-18-56	R-26 S1	REG	SW-846:9060	0	1

Only results shown in Section 13 'Display Flagged Data' are current as of this report generation. All other sections are valid for the date the COC data was inserted into EIM, and may have changed due to data updates in the intervening time.

March 16, 2018

Ms. Nita Patel
Los Alamos National Laboratory
TA-00, SM1237, Rm104C
Los Alamos, New Mexico 87545


Re: LANL- WQH Water Samples
Work Order: 444566
SDG: 2018-1811

Dear Ms. Patel:

GEL Laboratories, LLC (GEL) appreciates the opportunity to provide the following analytical results for the sample(s) we received on February 23, 2018, and analyzed for Explosives by LCMSMS, GC/MS Semivolatile, GC/MS Volatile, General Chemistry, Metals, Perchlorates by LCMSMS and Radiochemistry. This original data report has been prepared and reviewed in accordance with GEL's standard operating procedures.

Our policy is to provide high quality, personalized analytical services to enable you to meet your analytical needs on time every time. We trust that you will find everything in order and to your satisfaction. If you have any questions, please do not hesitate to call me at (843) 556-8171, ext. 4485.

Sincerely,


Brielle Luthman for
Valerie Davis
Project Manager

Chain of Custody: 2018-1811
Enclosures



ARS International, LLC (ARS-LANS-MTOA6-25093-GEL)
LANL- WQH Water Samples
Work Order #: 444566
SDG: 2018-1811

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Case Narrative

**Case Narrative for
ARS International, LLC (ARS-LANS-MTOA6-25093-GEL)
LANL- WQH Water Samples
Workorder #: 444566
SDG # : 2018-1811**

March 16, 2018

Laboratory Identification:

GEL Laboratories LLC
2040 Savage Road
Charleston, South Carolina 29407
(843) 556-8171

Summary

Sample receipt The samples arrived at GEL Laboratories LLC, Charleston, South Carolina on February 23, 2018 for analysis. The samples were delivered with proper chain of custody documentation and signatures. The samples were screened according to GEL Standard Operating Procedure. All sample containers arrived without any visible signs of tampering or breakage. Containers were checked for pH, where appropriate, and matched the preservative as documented on the accompanying chain of custody. Shipping container temperatures were checked, documented, and within specifications. Shipping container temperature was within specification (0 - 6C). There are no additional comments concerning sample receipt.

Sample Identification The laboratory received the following samples:

<u>Laboratory ID</u>	<u>Client ID</u>
444566001	CAPA-18-1
444566002	CAPA-18-2
444566003	CAPA-18-2
444566004	CAPA-18-6
444566005	CAPA-18-10
444566006	CAPA-18-12
444566007	CAPA-18-12
444566008	CAWA-18-54
444566009	CAWA-18-54
444566010	CAWA-18-111
444566011	CAWA-18-55
444566012	CAWA-18-56
444566013	CAWA-18-56
444566014	CAWA-18-112

Case Narrative


Sample analyses were conducted using methodology as outlined in GEL Laboratories, LLC (GEL) Standard Operating Procedures. Any technical or administrative problems during analysis, data review, and reduction are contained in the analytical case narratives in the enclosed data package.

Data Package

The enclosed data package contains the following sections: Case Narrative, Chain of Custody, Cooler Receipt

Checklist, Data Package Qualifier Definitions and data from the following fractions: Explosives by LCMSMS, GC/MS Semivolatile, GC/MS Volatile, General Chemistry, Metals, Perchlorates by LCMSMS and Radiochemistry.

I certify that this data report is in compliance with the terms and conditions of the subcontract and task order, both technically and for completeness, for other than the conditions detailed in the attached case narrative.


Brielle Luthman for
Valerie Davis
Project Manager

List of current GEL Certifications as of 16 March 2018

State	Certification
Alaska	17-018
Arkansas	88-0651
CLIA	42D0904046
California	2940
Colorado	SC00012
Connecticut	PH-0169
Delaware	SC00012
DoD ELAP/ ISO17025 A2LA	2567.01
Florida NELAP	E87156
Foreign Soils Permit	P330-15-00283, P330-15-00253
Georgia	SC00012
Georgia SDWA	967
Hawaii	SC00012
Idaho Chemistry	SC00012
Idaho Radiochemistry	SC00012
Illinois NELAP	200029
Indiana	C-SC-01
Kansas NELAP	E-10332
Kentucky SDWA	90129
Kentucky Wastewater	90129
Louisiana NELAP	03046 (AI33904)
Louisiana SDWA	LA180011
Maryland	270
Massachusetts	M-SC012
Michigan	9976
Mississippi	SC00012
Nebraska	NE-OS-26-13
Nevada	SC000122018-1
New Hampshire NELAP	205415
New Jersey NELAP	SC002
New Mexico	SC00012
New York NELAP	11501
North Carolina	233
North Carolina SDWA	45709
North Dakota	R-158
Oklahoma	9904
Pennsylvania NELAP	68-00485
Puerto Rico	SC00012
S. Carolina Radiochem	10120002
South Carolina Chemistry	10120001
Tennessee	TN 02934
Texas NELAP	T104704235-18-13
Utah NELAP	SC000122017-25
Vermont	VT87156
Virginia NELAP	460202
Washington	C780
West Virginia	997404

Chain of Custody and Supporting Documentation



Laboratories LLC

SAMPLE RECEIPT & REVIEW FORM

Client: <u>ESHL</u>		SDG/AR/COC/Work Order: <u>4445706</u>	
Received By: <u>ZKW</u>		Date Received: <u>2/23/18</u>	
Carrier and Tracking Number		Circle Applicable: <input checked="" type="checkbox"/> FedEx Express <input type="checkbox"/> FedEx Ground <input type="checkbox"/> UPS <input type="checkbox"/> Field Services <input type="checkbox"/> Courier <input type="checkbox"/> Other <u>5908 1783 5418-20c</u> <u>5908 1783 5430-2c</u> <u>5908 1783 5440-3c</u> <u>5908 1783 5429-2c</u>	
Suspected Hazard Information	Yes <input type="checkbox"/> No <input checked="" type="checkbox"/>	*If Net Counts > 100cpm on samples not marked "radioactive", contact the Radiation Safety Group for further investigation.	
Shipped as a DOT Hazardous?	Yes <input type="checkbox"/> No <input checked="" type="checkbox"/>	Hazard Class Shipped: _____ UN#: _____	
COC/Samples marked or classified as radioactive?	Yes <input type="checkbox"/> No <input checked="" type="checkbox"/>	Maximum Net Counts Observed* (Observed Counts - Area Background Counts): <u>0</u> <u>CPM</u> mR/Hr Classified as: Rad 1 Rad 2 Rad 3	
Is package, COC, and/or Samples marked HAZ?	Yes <input checked="" type="checkbox"/> No <input type="checkbox"/>	If yes, select Hazards below, and contact the GEL Safety Group. <input checked="" type="checkbox"/> PCB's <input type="checkbox"/> Flammable <input type="checkbox"/> Foreign Soil <input type="checkbox"/> RCRA <input type="checkbox"/> Asbestos <input type="checkbox"/> Beryllium <input type="checkbox"/> Other: _____	

Sample Receipt Criteria	Yes	NA	No	Comments/Qualifiers (Required for Non-Conforming Items)
1 Shipping containers received intact and sealed?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	Circle Applicable: Seals broken Damaged container Leaking container Other (describe)
2 Chain of custody documents included with shipment?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
3 Samples requiring cold preservation within (0 ≤ 6 deg. C)?*	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	Preservation Method: Wet Ice <input checked="" type="checkbox"/> Ice Packs Dry ice <input type="checkbox"/> <u>None</u> Other: _____ *all temperatures are recorded in Celsius TEMP: <u>See Above</u>
4 Daily check performed and passed on IR temperature gun?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	Temperature Device Serial #: <u>IR3-16</u> Secondary Temperature Device Serial # (If Applicable): _____
5 Sample containers intact and sealed?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	Circle Applicable: Seals broken Damaged container Leaking container Other (describe)
6 Samples requiring chemical preservation at proper pH?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	Sample ID's and Containers Affected: _____ If Preservation added, Lot#: _____
7 Do any samples require Volatile Analysis?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	If Yes, Are Encores or Soil Kits present? Yes _____ No <input checked="" type="checkbox"/> (If yes, take to VOA Freezer) Do VOA vials contain acid preservation? Yes _____ No <input checked="" type="checkbox"/> N/A _____ (If unknown, select No) VOA vials free of headspace? Yes <input checked="" type="checkbox"/> No _____ N/A _____ Sample ID's and containers affected: _____
8 Samples received within holding time?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	ID's and tests affected: _____
9 Sample ID's on COC match ID's on bottles?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	Sample ID's and containers affected: _____
10 Date & time on COC match date & time on bottles?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	Sample ID's affected: _____
11 Number of containers received match number indicated on COC? <u>2/23/18</u>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	Sample ID's affected: <u>We only rec'd 2 HEXMOD Cont. for -18-12</u>
12 Are sample containers identifiable as GEL provided?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
13 COC form is properly signed in relinquished/received sections?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	

Comments (Use Continuation Form if needed):

PM (or PMA) review: Initials BL Date 2/23/18 Page 1 of 1

GL-CHL-SR-001 Rev 5

ORIGIN ID:SAFA (505) 665-9966
KEITH GREENE
LOS ALAMOS NATL LAB.
TA00 BLDG 1237 DPU 03

SHIP DATE: 22FEB18
ACTWGT: 48.0 LB MAN
CAD: 0014176/CAFE2916

LOS ALAMOS, NM 87545
UNITED STATES US

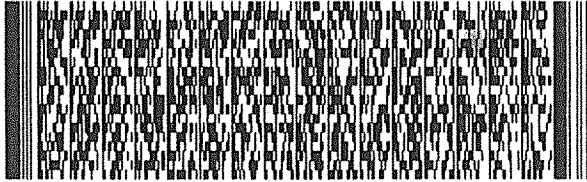
BILL SENDER

TO **VALERIE DAVIS**
GENERAL ENGINEERING LAB
2040 SAVAGE RD

CHARLESTON SC 29407

(843) 666-8171

REF: 8A030AWE6L11551000



FedEx
Express



J1513150813011v

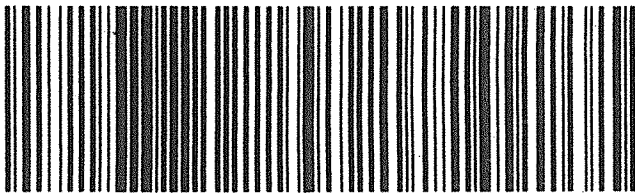
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FRI - 23 FEB 10:30A
PRIORITY OVERNIGHT

0201

X7 RBWA

29407
SC-US CHS



ORIGIN ID:SAFA (505) 665-9966
KEITH GREENE
LOS ALAMOS NATL LAB.
TA00 BLDG 1237 DPU 03

SHIP DATE: 22FEB18
ACTWGT: 46.0 LB MAN
CAD: 0014176/CAFE2916

LOS ALAMOS, NM 87545
UNITED STATES US

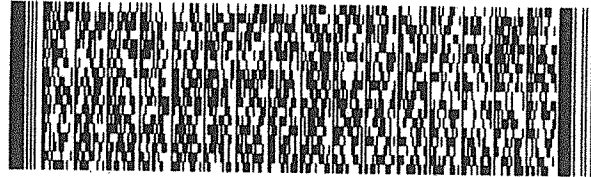
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TO **VALERIE DAVIS**
GENERAL ENGINEERING LAB
2040 SAVAGE RD

CHARLESTON SC 29407

(843) 666-8171

REF: 21PD0ASRD06BDBD00



FedEx
Express



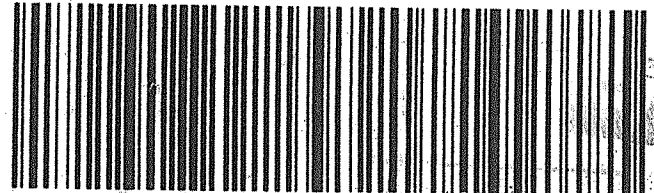
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PRIORITY OVERNIGHT

X7 RBWA

29407
SC-US CHS



ORIGIN ID:SAFA (505) 665-9966
KEITH GREENE
LOS ALAMOS NATL LAB.
TA00 BLDG 1237 DPU 03

SHIP DAT
ACTWGT:
CAD: 00

LOS ALAMOS, NM 87545
UNITED STATES US

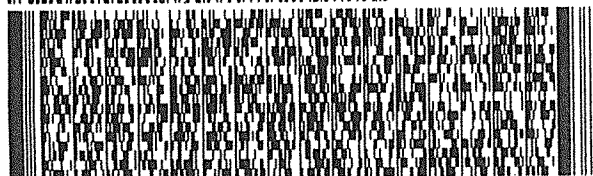
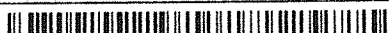
BILL SE

TO **VALERIE DAVIS**
GENERAL ENGINEERING LAB
2040 SAVAGE RD

CHARLESTON SC 29407

(843) 556-8171

REF: 8A030AWE6L11551000



FedEx
Express



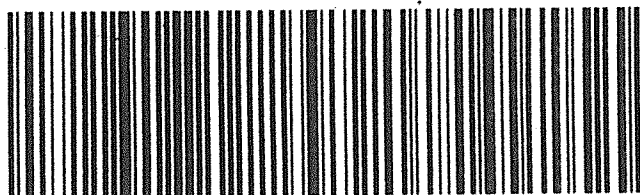
1 of 3

TRK# 5908 1783 5418
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MASTER

FRI - 23 FEB 10:30A
PRIORITY OVERNIGHT

X7 RBWA

29407
SC-US CHS



Part # 156148V-434 RIT2 EXP 02/18 ***

ORIGIN ID:SAFA (505) 665-9966
KEITH GREENE
LOS ALAMOS NATL LAB.
TA00 BLDG 1237 DPU 03

SHIP DATE: 22FEB18
ACTWGT: 50.0 LB MAN
CAD: 0014176/CAFE2916

LOS ALAMOS, NM 87545
UNITED STATES US

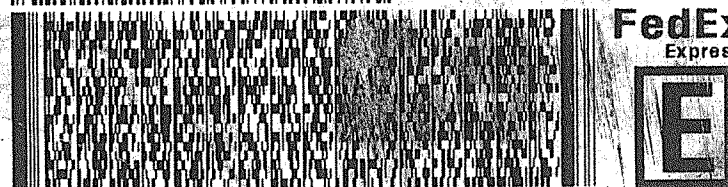
BILL SENDER

TO **VALERIE DAVIS**
GENERAL ENGINEERING LAB
2040 SAVAGE RD

CHARLESTON SC 29407

(843) 556-8171

REF: 8A030AWE6L11551000



FedEx
Express



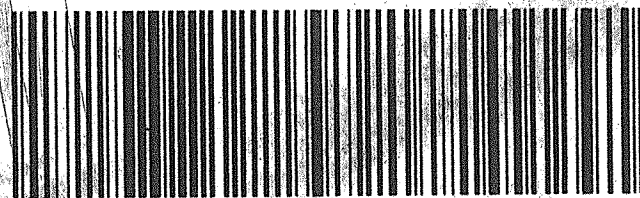
3 of 3

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0263
Mstr# 5908 1783 5418

FRI - 23 FEB 10:30
PRIORITY OVERNIGHT

X7 RBWA

2940
SC-US CH



Part # 156148V-434 RIT2 EXP 02/18 ***

Data Review Qualifier Flag Definition Sheet

Data Review Qualifier Definitions

Qualifier	Explanation
-----------	-------------

*	A quality control analyte recovery is outside of specified acceptance criteria
**	Analyte is a surrogate compound
<	Result is less than value reported
>	Result is greater than value reported
^	RPD of sample and duplicate evaluated using +/-RL. Concentrations are <5X the RL
A	The TIC is a suspected aldol-condensation product
B	Target analyte was detected in the associated blank
B	Metals-Either presence of analyte detected in the associated blank, or MDL/IDL < sample value < PQL
BD	Results are either below the MDC or tracer recovery is low
C	Analyte has been confirmed by GC/MS analysis
D	Results are reported from a diluted aliquot of the sample
d	5-day BOD-The 2:1 depletion requirement was not met for this sample
E	Organics-Concentration of the target analyte exceeds the instrument calibration range
E	Metals-%difference of sample and SD is >10%. Sample concentration must meet flagging criteria
H	Analytical holding time was exceeded
h	Preparation or preservation holding time was exceeded
J	Value is estimated
N	Metals-The Matrix spike sample recovery is not within specified control limits
N	Organics-Presumptive evidence based on mass spectral library search to make a tentative identification of the analyte (TIC). Quantitation is based on nearest internal standard response factor
N/A	Spike recovery limits do not apply. Sample concentration exceeds spike concentration by 4X or more
ND	Analyte concentration is not detected above the reporting limit
UI	Gamma Spectroscopy-Uncertain identification
X	Consult Case Narrative, Data Summary package, or Project Manager concerning this qualifier
Y	QC Samples were not spiked with this compound
Z	Paint Filter Test-Particulates passed through the filter, however no free liquids were observed.

P Organics-The concentrations between the primary and confirmation columns/detectors is >40% difference.
For HPLC, the difference is >70%.

U Analyte was analyzed for, but not detected above the MDL, MDA, or LOD.

Volatile Analysis

Case Narrative

**GC/MS Volatile
Technical Case Narrative
ARS International, LLC (ARSL)
SDG #: 2018-1811
Work Order #: 444566**

Method/Analysis Information

Procedure: Volatile Organic Compounds (VOC) by Gas Chromatograph/Mass Spectrometer

Analytical Method: SW-846:8260B

Analytical Batch
Number: 1744716

Sample Analysis

The following client and quality control samples were analyzed to complete this SDG using the methods referenced in the Analysis Information section:

Sample ID	Client ID
444566002	CAPA-18-2
444566004	CAPA-18-6
444566006	CAPA-18-12
444566009	CAWA-18-54
444566010	CAWA-18-111
444566012	CAWA-18-56
444566014	CAWA-18-112
1203984431	Method Blank (MB)
1203984433	Laboratory Control Sample (LCS)
1203984434	Laboratory Control Sample (LCS)
1203984437	444617006(CAWA-18-67) Post Spike (PS)
1203984438	444617006(CAWA-18-67) Post Spike (PS)
1203984439	444617006(CAWA-18-67) Post Spike Duplicate (PSD)
1203984440	444617006(CAWA-18-67) Post Spike Duplicate (PSD)
1203985580	Method Blank (MB)
1203985581	Laboratory Control Sample (LCS)
1203985582	Laboratory Control Sample (LCS)

NOTE: For volatile organic analyses the matrix spike designations may be indicated as "PS" or "PSD". The "PS" designation (post spike) indicates that the matrix was fortified prior to analysis but after applying any prep factors, such as a dilution. The laboratory considers the MS/MSD and PS/PSD designations interchangeable.

The data results reported met all SOP and method criteria, unless otherwise discussed below.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-OA-E-038 REV# 26.

Calibration Information

A complete list of the initial calibration data files with the correct dates and times of analysis are shown in the Calibration History report located in the Standard Data section of the data package. The surrogate compounds were calibrated using a minimum five-point calibration curve. The surrogates were added by the auto sampler at a concentration of 50 ug/L or 20 ug/L for low level analyses. GEL Laboratories LLC will not have surrogate recoveries reported for Dibromofluoromethane. This is due to increased regulations for this analyte and an industry shortage.

Initial Calibration

All initial calibration requirements have been met for this sample delivery group (SDG).

Continuing Calibration Verification Requirements

All associated calibration verification standard(s) (CCV) met the acceptance criteria.

Quality Control (QC) Information

Blank (MB) Statement

The blanks analyzed with this SDG met the acceptance criteria.

Surrogate Recoveries

Surrogate recoveries in all client and quality control samples were within the acceptance limits.

Laboratory Control Sample (LCS) Recovery

The LCS/and or LCSD (See Below) recoveries were not all within the acceptance limits. The unacceptable recoveries were less than 5% of the requested analyte list. This satisfies the client criteria. The results are reported.

Sample	Analyte	Value
1203985581 (LCS)	Hexachlorobutadiene	66* (72%-136%)

QC Sample Designation

Sample 444617006 (CAWA-18-67) was designated for spike analysis.

Matrix Spike/Matrix Spike Duplicate Recovery Statement

The matrix spike (MS) and matrix spike duplicate (MSD) recoveries were within the required acceptance limits.

Relative Percent Difference (RPD) Statement

The RPDs between the matrix spike pair met the acceptance limits.

Internal Standard (ISTD) Acceptance

The internal standard responses in all client and quality control samples met the required acceptance criteria.

Technical Information

Holding Time Specifications

All samples in this SDG met the specified holding time. GEL assigns holding times based on the associated methodology, which assigns the date and time from sample collection or sample receipt. Those holding times expressed in hours are calculated in the ALPHALIMS system. Those holding times expressed as days expire at midnight on the day of expiration.

Sample Preservation and Integrity

All samples met the sample preservation and integrity requirements.

Sample Dilutions/Methanol Dilutions

The samples in this SDG did not require dilutions.

Sample Re-extraction/Re-analysis

Re-analyses were not required for samples in this SDG.

Miscellaneous Information**Manual Integrations**

Data files associated with the initial calibration, continuing calibration check, and samples did not require manual integrations.

TIC Comment

Tentatively identified compounds (TIC) may be requested for samples 444566002 (CAPA-18-2), 444566004 (CAPA-18-6), 444566006 (CAPA-18-12), 444566009 (CAWA-18-54), 444566010 (CAWA-18-111), 444566012 (CAWA-18-56) and 444566014 (CAWA-18-112) in this delivery group/work order. Please note that non-requested calibrated analytes detected in a client sample may be reported on the Form 1/Certificate of Analysis as TICs. TIC data, if requested, were included on the Sample Data Summary (Form 1) and included with the sample raw data.

Additional Comments

Additional comments were not required for this SDG.

Residual Chlorine

Residual Chlorine was not detected in any of the samples in this SDG.

Electronic Package Comment

The following package was generated using an electronic data processing program referred to as "virtual packaging". In an effort to increase quality and efficiency, the laboratory is developing systems to eventually generate all data packages electronically. The following change from "traditional" packages should be noted: Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. An electronic signature page inserted after the case narrative of each electronic package will indicate the reviewer name associated with the generation of the data and package. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

System Configuration

The Volatile-GC/MS analysis was performed on the following instrument configuration:

Instrument ID	Instrument	System Configuration	Column ID	Column Description	P & T Trap
VOA9.I	Agilent 6890/5973 GC/MS w/ OI Eclipse/Archon Autosampler	HP6890/HP5973	DB-624	J&W, 60m x 0.25mm x 1.4um	Trap 10

Certification Statement

Where the analytical method has been performed under NELAP certification, the analysis has met all of the requirements of the NELAC standard unless otherwise noted in the analytical case narrative.

GEL LABORATORIES LLC

2040 Savage Road Charleston SC 29407 - (843) 556-8171 - www.gel.com

Qualifier Definition Report for

ARSL004 ARS International, LLC (ARS-LANS-MTOA6-25093-GEL)

Client SDG: 2018-1811 GEL Work Order: 444566

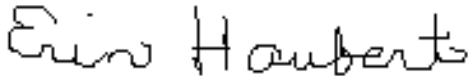
The Qualifiers in this report are defined as follows:

- * A quality control analyte recovery is outside of specified acceptance criteria
- ** Analyte is a surrogate compound
- J Value is estimated
- U Analyte was analyzed for, but not detected above the MDL, MDA, MDC or LOD.
- DL Indicates that sample is diluted.
- RA Indicates that sample is re-analyzed without re-extraction.
- RE Indicates that sample is re-extracted.

Review/Validation

GEL requires all analytical data to be verified by a qualified data reviewer. In addition, all CLP-like deliverables receive a third level review of the fractional data package.

The following data validator verified the information presented in this data report:

Signature: 

Name: Erin Haubert

Date: 21 MAR 2018

Title: Data Validator

Sample Data Summary

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566002

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 03:36

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 03:36

Data File: 030618V9\91234.D

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	J	2.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

Page 2 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566002

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 03:36

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 03:36

Data File: 030618V9\9I234.D

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811

Lab Sample ID: 444566002

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Client ID: CAPA-18-2

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 03:36

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 03:36

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	52.2	50.0	ug/L 104	(71%-134%)
Bromofluorobenzene	49.5	50.0	ug/L 99	(70%-131%)
Toluene-d8	51.0	50.0	ug/L 102	(74%-124%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
	unknown siloxane	14.651	20.5	ug/L	0	J
	unknown siloxane	16.62	6.23	ug/L	0	J

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566004

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Batch ID: 1744716

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Run Date: 03/07/2018 02:09

Inst: VOA9.I

Dilution: 1

Prep Date: 03/07/2018 02:09

Analyst: RXY1

Purge Vol: 5 mL

Data File: 030618V9\91231.D

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

Page 2 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566004

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 02:09

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 02:09

Data File: 030618V9\9I231.D

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

Page 3 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566004

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Client ID: CAPA-18-6

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 02:09

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 02:09

Column: DB-624

Data File: 030618V9\9I231.D

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	51.2	50.0	ug/L 102	(71%-134%)
Bromofluorobenzene	50.5	50.0	ug/L 101	(70%-131%)
Toluene-d8	52.1	50.0	ug/L 104	(74%-124%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
	unknown siloxane	14.651	24.6	ug/L	0	J

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566006

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Client ID: CAPA-18-12

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 04:04

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 04:04

Column: DB-624

Data File: 030618V9\91235.D

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

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SDG Number: 2018-1811

Lab Sample ID: 444566006

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 04:04

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 04:04

Data File: 030618V9\9I235.D

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811

Lab Sample ID: 444566006

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Client ID: CAPA-18-12

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 04:04

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 04:04

Column: DB-624

Data File: 030618V9\9I235.D

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	52.5	50.0	ug/L 105	(71%-134%)
Bromofluorobenzene	51.5	50.0	ug/L 103	(70%-131%)
Toluene-d8	54.6	50.0	ug/L 109	(74%-124%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
	unknown siloxane	12.291	6.13	ug/L	0	J
	unknown siloxane	14.651	35	ug/L	0	J
	unknown siloxane	16.62	8.16	ug/L	0	J

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566009

Date Collected: 02/21/2018 10:31

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Client ID: CAWA-18-54

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 04:33

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 04:33

Column: DB-624

Data File: 030618V9\91236.D

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	J	1.62	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811

Lab Sample ID: 444566009

Date Collected: 02/21/2018 10:31

Date Received: 02/23/2018 09:25

Matrix: W

Client ID: CAWA-18-54

Batch ID: 1744716

Run Date: 03/07/2018 04:33

Prep Date: 03/07/2018 04:33

Data File: 030618V9\9I236.D

Client: ARSL004

Method: SW-846:8260B

Inst: VOA9.I

Analyst: RXY1

Project: ESHL00114

SOP Ref: GL-OA-E-038

Dilution: 1

Purge Vol: 5 mL

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene		1.41	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811

Lab Sample ID: 444566009

Date Collected: 02/21/2018 10:31

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 04:33

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 04:33

Data File: 030618V9\9I236.D

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	52.1	50.0	ug/L 104	(71%-134%)
Bromofluorobenzene	51.6	50.0	ug/L 103	(70%-131%)
Toluene-d8	52.6	50.0	ug/L 105	(74%-124%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
	unknown siloxane	14.651	27.7	ug/L	0	J
	unknown siloxane	16.62	5.7	ug/L	0	J

Volatile
Certificate of Analysis
Sample Summary

SDG Number: 2018-1811

Lab Sample ID: 444566010

Date Collected: 02/21/2018 10:31

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 02:38

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 02:38

Data File: 030618V9\91232.D

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

Page 2 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566010

Date Collected: 02/21/2018 10:31

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 02:38

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 02:38

Data File: 030618V9\91232.D

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

Page 3 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566010

Date Collected: 02/21/2018 10:31

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 02:38

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 02:38

Column: DB-624

Data File: 030618V9\9I232.D

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	51.0	50.0	ug/L 102	(71%-134%)
Bromofluorobenzene	48.8	50.0	ug/L 98	(70%-131%)
Toluene-d8	51.4	50.0	ug/L 103	(74%-124%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
	unknown siloxane	14.651	21.4	ug/L	0	J

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566012

Date Collected: 02/21/2018 10:44

Date Received: 02/23/2018 09:25

Matrix: W

Client ID: CAWA-18-56

Batch ID: 1744716

Run Date: 03/07/2018 05:01

Prep Date: 03/07/2018 05:01

Data File: 030618V9\91237.D

Client: ARSL004

Method: SW-846:8260B

Inst: VOA9.I

Analyst: RXY1

Project: ESHL00114

SOP Ref: GL-OA-E-038

Dilution: 1

Purge Vol: 5 mL

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

Page 2 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566012

Date Collected: 02/21/2018 10:44

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 05:01

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 05:01

Data File: 030618V9\91237.D

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811

Lab Sample ID: 444566012

Date Collected: 02/21/2018 10:44

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Client ID: CAWA-18-56

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 05:01

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 05:01

Column: DB-624

Data File: 030618V9\9I237.D

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	53.5	50.0	ug/L 107	(71%-134%)
Bromofluorobenzene	53.0	50.0	ug/L 106	(70%-131%)
Toluene-d8	52.2	50.0	ug/L 104	(74%-124%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
	unknown siloxane	14.651	28.9	ug/L	0	J
	unknown siloxane	16.62	6.45	ug/L	0	J

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566014

Date Collected: 02/21/2018 10:44

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Client ID: CAWA-18-112

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 03:07

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 03:07

Column: DB-624

Data File: 030618V9\91233.D

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

SDG Number: 2018-1811

Lab Sample ID: 444566014

Date Collected: 02/21/2018 10:44

Date Received: 02/23/2018 09:25

Matrix: W

Client ID: CAWA-18-112

Batch ID: 1744716

Run Date: 03/07/2018 03:07

Prep Date: 03/07/2018 03:07

Data File: 030618V9\9I233.D

Client: ARSL004

Method: SW-846:8260B

Inst: VOA9.I

Analyst: RXY1

Project: ESHL00114

SOP Ref: GL-OA-E-038

Dilution: 1

Purge Vol: 5 mL

Column: DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

Page 3 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566014

Date Collected: 02/21/2018 10:44

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Client ID: CAWA-18-112

Method: SW-846:8260B

SOP Ref: GL-OA-E-038

Batch ID: 1744716

Inst: VOA9.I

Dilution: 1

Run Date: 03/07/2018 03:07

Analyst: RXY1

Purge Vol: 5 mL

Prep Date: 03/07/2018 03:07

Column: DB-624

Data File: 030618V9\9I233.D

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	50.8	50.0	ug/L 102	(71%-134%)
Bromofluorobenzene	49.2	50.0	ug/L 98	(70%-131%)
Toluene-d8	51.1	50.0	ug/L 102	(74%-124%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
	unknown siloxane	12.291	5.44	ug/L	0	J
	unknown siloxane	14.651	27.6	ug/L	0	J
	unknown siloxane	16.62	5.27	ug/L	0	J

Quality Control Summary

Volatile
Surrogate Recovery Report

Page 1 of 1

SDG Number: 2018-1811**Matrix Type: LIQUID**

Sample ID	Client ID	DCED4 %REC	TOL %REC	BFB %REC
1203984433	LCS for batch 1744716	99	103	99
1203984434	LCS for batch 1744716	100	106	103
1203984431	MB for batch 1744716	98	101	99
444566004	CAPA-18-6	102	104	101
444566010	CAWA-18-111	102	103	98
444566014	CAWA-18-112	102	102	98
444566002	CAPA-18-2	104	102	99
444566006	CAPA-18-12	105	109	103
444566009	CAWA-18-54	104	105	103
444566012	CAWA-18-56	107	104	106
1203985581	LCS for batch 1744716	97	99	99
1203985582	LCS for batch 1744716	100	103	97
1203985580	MB for batch 1744716	99	103	100
1203984437	CAWA-18-67PS	100	103	103
1203984439	CAWA-18-67PSD	103	107	104
1203984438	CAWA-18-67PS	97	102	97
1203984440	CAWA-18-67PSD	98	103	101

Surrogate**Acceptance Limits**

DCED4 = 1,2-Dichloroethane-d4

(71%-134%)

TOL = Toluene-d8

(74%-124%)

BFB = Bromofluorobenzene

(70%-131%)

* Recovery outside Acceptance Limits

Column to be used to flag recovery values

D Sample Diluted

Volatile
Quality Control Summary
Spike Recovery Report

Page 1 of 4

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203984433

Instrument: VOA9.I

Analysis Date: 03/07/2018 00:14

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
179601-23-1	LCS m,p-Xylenes	100	0.0	83.9	84	71-127
75-05-8	LCS Acetonitrile	1250	0.0	1120	90	61-125
67-64-1	LCS Acetone	250	0.0	198	79	48-157
74-88-4	LCS Iodomethane	250	0.0	201	80	72-128
75-15-0	LCS Carbon disulfide	250	0.0	193	77	69-138
108-05-4	LCS Vinyl acetate	250	0.0	233	93	67-125
78-93-3	LCS 2-Butanone	250	0.0	215	86	55-138
108-10-1	LCS 4-Methyl-2-pentanone	250	0.0	233	93	66-124
591-78-6	LCS 2-Hexanone	250	0.0	212	85	56-140
75-71-8	LCS Dichlorodifluoromethane	50.0	0.0	42.2	84	40-160
74-87-3	LCS Chloromethane	50.0	0.0	46.8	94	58-135
75-01-4	LCS Vinyl chloride	50.0	0.0	47.9	96	65-137
74-83-9	LCS Bromomethane	50.0	0.0	40.9	82	63-137
75-00-3	LCS Chloroethane	50.0	0.0	41.9	84	69-129
75-69-4	LCS Trichlorofluoromethane	50.0	0.0	41.5	83	69-138
60-29-7	LCS Ethyl ether	50.0	0.0	48.4	97	72-125
75-35-4	LCS 1,1-Dichloroethylene	50.0	0.0	38.8	78	66-126
75-09-2	LCS Methylene chloride	50.0	0.0	39.9	80	68-119
1634-04-4	LCS tert-Butyl methyl ether	50.0	0.0	45.6	91	76-128
156-60-5	LCS trans-1,2-Dichloroethylene	50.0	0.0	40.1	80	71-124
75-34-3	LCS 1,1-Dichloroethane	50.0	0.0	41.5	83	73-123
156-59-2	LCS cis-1,2-Dichloroethylene	50.0	0.0	41.3	83	75-123

Volatile
Quality Control Summary
Spike Recovery Report

Page 2 of 4

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203984433

Instrument: VOA9.I

Analysis Date: 03/07/2018 00:14

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
594-20-7	LCS 2,2-Dichloropropane	50.0	0.0	38.5	77	72-138
74-97-5	LCS Bromochloromethane	50.0	0.0	44.9	90	76-125
67-66-3	LCS Chloroform	50.0	0.0	41.5	83	76-123
71-55-6	LCS 1,1,1-Trichloroethane	50.0	0.0	40.0	80	74-136
563-58-6	LCS 1,1-Dichloropropene	50.0	0.0	37.5	75	72-129
56-23-5	LCS Carbon tetrachloride	50.0	0.0	40.4	81	72-140
107-06-2	LCS 1,2-Dichloroethane	50.0	0.0	42.8	86	74-122
71-43-2	LCS Benzene	50.0	0.0	40.1	80	72-121
79-01-6	LCS Trichloroethylene	50.0	0.0	40.1	80	74-125
78-87-5	LCS 1,2-Dichloropropane	50.0	0.0	41.7	83	73-121
74-95-3	LCS Dibromomethane	50.0	0.0	44.5	89	78-123
75-27-4	LCS Bromodichloromethane	50.0	0.0	45.6	91	77-131
10061-01-5	LCS cis-1,3-Dichloropropylene	50.0	0.0	44.9	90	78-131
108-88-3	LCS Toluene	50.0	0.0	40.8	82	71-121
10061-02-6	LCS trans-1,3-Dichloropropylene	50.0	0.0	48.2	96	78-131
79-00-5	LCS 1,1,2-Trichloroethane	50.0	0.0	46.0	92	74-118
142-28-9	LCS 1,3-Dichloropropane	50.0	0.0	44.1	88	74-118
127-18-4	LCS Tetrachloroethylene	50.0	0.0	39.4	79	69-129
124-48-1	LCS Dibromochloromethane	50.0	0.0	45.6	91	76-137
106-93-4	LCS 1,2-Dibromoethane	50.0	0.0	47.3	95	78-122
108-90-7	LCS Chlorobenzene	50.0	0.0	41.7	83	74-120
100-41-4	LCS Ethylbenzene	50.0	0.0	40.3	81	73-125

Volatile
Quality Control Summary
Spike Recovery Report

Page 3 of 4

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203984433

Instrument: VOA9.I

Analysis Date: 03/07/2018 00:14

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
95-47-6	LCS o-Xylene	50.0	0.0	42.1	84	74-126
100-42-5	LCS Styrene	50.0	0.0	44.8	90	72-130
75-25-2	LCS Bromoform	50.0	0.0	55.7	111	72-136
98-82-8	LCS Isopropylbenzene	50.0	0.0	41.5	83	70-130
79-34-5	LCS 1,1,2,2-Tetrachloroethane	50.0	0.0	48.7	97	70-126
96-18-4	LCS 1,2,3-Trichloropropane	50.0	0.0	49.3	99	74-122
108-86-1	LCS Bromobenzene	50.0	0.0	43.3	87	74-120
103-65-1	LCS n-Propylbenzene	50.0	0.0	39.7	79	67-128
108-67-8	LCS 1,3,5-Trimethylbenzene	50.0	0.0	42.2	84	70-129
95-49-8	LCS 2-Chlorotoluene	50.0	0.0	42.0	84	71-124
106-43-4	LCS 4-Chlorotoluene	50.0	0.0	40.5	81	69-125
98-06-6	LCS tert-Butylbenzene	50.0	0.0	41.8	84	72-130
95-63-6	LCS 1,2,4-Trimethylbenzene	50.0	0.0	41.8	84	70-126
135-98-8	LCS sec-Butylbenzene	50.0	0.0	40.7	81	70-131
99-87-6	LCS 4-Isopropyltoluene	50.0	0.0	41.0	82	71-131
541-73-1	LCS 1,3-Dichlorobenzene	50.0	0.0	41.2	82	72-121
106-46-7	LCS 1,4-Dichlorobenzene	50.0	0.0	41.5	83	71-120
104-51-8	LCS n-Butylbenzene	50.0	0.0	38.4	77	68-134
96-12-8	LCS 1,2-Dibromo-3-chloropropane	50.0	0.0	47.9	96	68-141
87-68-3	LCS Hexachlorobutadiene	50.0	0.0	40.7	81	72-136
91-20-3	LCS Naphthalene	50.0	0.0	52.0	104	72-132
87-61-6	LCS 1,2,3-Trichlorobenzene	50.0	0.0	47.3	95	70-130

Volatile

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Quality Control Summary
Spike Recovery Report

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203984433

Instrument: VOA9.I

Analysis Date: 03/07/2018 00:14

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
120-82-1	LCS 1,2,4-Trichlorobenzene	50.0	0.0	44.2	88	71-129
630-20-6	LCS 1,1,1,2-Tetrachloroethane	50.0	0.0	46.5	93	79-127
95-50-1	LCS 1,2-Dichlorobenzene	50.0	0.0	42.9	86	74-120
71-36-3	LCS n-Butyl alcohol	5000	0.0	5020	100	63-138

Volatile
Quality Control Summary
Spike Recovery Report

Page 1 of 1

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203984434

Instrument: VOA9.I

Analysis Date: 03/07/2018 01:11

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
107-02-8	LCS Acrolein	250	0.0	251	101	60-140
76-13-1	LCS Trichlorotrifluoroethane	250	0.0	227	91	61-148
107-05-1	LCS Allyl chloride	250	0.0	222	89	59-125
107-13-1	LCS Acrylonitrile	250	0.0	259	104	65-122
107-12-0	LCS Propionitrile	250	0.0	255	102	64-124
126-98-7	LCS Methacrylonitrile	250	0.0	248	99	64-126
80-62-6	LCS Methyl methacrylate	250	0.0	251	100	69-127
97-63-2	LCS Ethyl methacrylate	250	0.0	242	97	66-130
78-83-1	LCS Isobutyl alcohol	2500	0.0	2630	105	65-135
126-99-8	LCS 2-Chloro-1,3-butadiene	50.0	0.0	47.4	95	66-147

Volatile
Quality Control Summary
Spike Recovery Report

Page 1 of 8

SDG Number: 2018-1811

Sample Type: Post Spike

Client ID: CAWA-18-67PS

Matrix: W

Lab Sample ID 1203984437

Instrument: VOA9.I

Analysis Date: 03/08/2018 02:43

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
179601-23-1	PS m,p-Xylenes	100	0.00 U	97.6	98	59-132
75-05-8	PS Acetonitrile	1250	0.00 U	1050	84	56-131
67-64-1	PS Acetone	250	0.00 U	144	58	25-155
74-88-4	PS Iodomethane	250	0.00 U	228	91	66-133
75-15-0	PS Carbon disulfide	250	0.00 U	226	90	61-141
108-05-4	PS Vinyl acetate	250	0.00 U	209	84	48-133
78-93-3	PS 2-Butanone	250	0.00 U	181	73	25-143
108-10-1	PS 4-Methyl-2-pentanone	250	0.00 U	216	87	61-127
591-78-6	PS 2-Hexanone	250	0.00 U	194	78	33-138
75-71-8	PS Dichlorodifluoromethane	50.0	0.00 U	49.5	99	33-164
74-87-3	PS Chloromethane	50.0	0.00 U	59.6	119	53-139
75-01-4	PS Vinyl chloride	50.0	0.00 U	59.8	120	58-140
74-83-9	PS Bromomethane	50.0	0.00 U	50.1	100	59-146
75-00-3	PS Chloroethane	50.0	0.00 U	50.5	101	65-129
75-69-4	PS Trichlorofluoromethane	50.0	0.00 U	49.0	98	65-141
60-29-7	PS Ethyl ether	50.0	0.00 U	50.3	101	69-127
75-35-4	PS 1,1-Dichloroethylene	50.0	0.00 U	46.4	93	59-130
75-09-2	PS Methylene chloride	50.0	0.00 U	45.0	90	62-123
1634-04-4	PS tert-Butyl methyl ether	50.0	0.00 U	47.9	96	69-132
156-60-5	PS trans-1,2-Dichloroethylene	50.0	0.00 U	46.9	94	65-127
75-34-3	PS 1,1-Dichloroethane	50.0	0.00 U	47.6	95	67-127
156-59-2	PS cis-1,2-Dichloroethylene	50.0	0.00 U	47.2	94	69-127

Volatile
Quality Control Summary
Spike Recovery Report

Page 2 of 8

SDG Number: 2018-1811

Sample Type: Post Spike

Client ID: CAWA-18-67PS

Matrix: W

Lab Sample ID 1203984437

Instrument: VOA9.I

Analysis Date: 03/08/2018 02:43

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
594-20-7	PS 2,2-Dichloropropane	50.0	0.00	U 44.6	89	66-137
74-97-5	PS Bromochloromethane	50.0	0.00	U 48.2	96	71-130
67-66-3	PS Chloroform	50.0	0.00	U 46.6	93	71-129
71-55-6	PS 1,1,1-Trichloroethane	50.0	0.00	U 47.6	95	69-139
563-58-6	PS 1,1-Dichloropropene	50.0	0.00	U 44.1	88	67-130
56-23-5	PS Carbon tetrachloride	50.0	0.00	U 49.1	98	66-143
107-06-2	PS 1,2-Dichloroethane	50.0	0.00	U 46.6	93	69-130
71-43-2	PS Benzene	50.0	0.00	U 46.3	93	66-125
79-01-6	PS Trichloroethylene	50.0	0.00	U 46.9	94	65-131
78-87-5	PS 1,2-Dichloropropane	50.0	0.00	U 47.6	95	67-127
74-95-3	PS Dibromomethane	50.0	0.00	U 47.9	96	72-129
75-27-4	PS Bromodichloromethane	50.0	0.00	U 52.1	104	70-138
10061-01-5	PS cis-1,3-Dichloropropylene	50.0	0.00	U 50.4	101	70-134
108-88-3	PS Toluene	50.0	0.00	U 45.9	92	60-126
10061-02-6	PS trans-1,3-Dichloropropylene	50.0	0.00	U 51.4	103	69-135
79-00-5	PS 1,1,2-Trichloroethane	50.0	0.00	U 48.3	97	66-125
142-28-9	PS 1,3-Dichloropropane	50.0	0.00	U 47.4	95	67-124
127-18-4	PS Tetrachloroethylene	50.0	0.00	U 46.5	93	60-130
124-48-1	PS Dibromochloromethane	50.0	0.00	U 49.3	99	68-143
106-93-4	PS 1,2-Dibromoethane	50.0	0.00	U 50.6	101	71-127
108-90-7	PS Chlorobenzene	50.0	0.00	U 47.3	95	64-124
100-41-4	PS Ethylbenzene	50.0	0.00	U 46.7	93	61-130

Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Post Spike

Client ID: CAWA-18-67PS

Matrix: W

Lab Sample ID 1203984437

Instrument: VOA9.I

Analysis Date: 03/08/2018 02:43

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
95-47-6	PS o-Xylene	50.0	0.00 U	48.0	96	62-131
100-42-5	PS Styrene	50.0	0.00 U	50.5	101	59-135
75-25-2	PS Bromoform	50.0	0.00 U	59.4	119	64-138
98-82-8	PS Isopropylbenzene	50.0	0.00 U	48.6	97	55-133
79-34-5	PS 1,1,2,2-Tetrachloroethane	50.0	0.00 U	50.3	101	62-129
96-18-4	PS 1,2,3-Trichloropropane	50.0	0.00 U	50.0	100	70-124
108-86-1	PS Bromobenzene	50.0	0.00 U	49.3	99	62-124
103-65-1	PS n-Propylbenzene	50.0	0.00 U	47.4	95	50-133
108-67-8	PS 1,3,5-Trimethylbenzene	50.0	0.00 U	49.4	99	53-135
95-49-8	PS 2-Chlorotoluene	50.0	0.00 U	49.4	99	56-128
106-43-4	PS 4-Chlorotoluene	50.0	0.00 U	47.2	94	53-130
98-06-6	PS tert-Butylbenzene	50.0	0.00 U	50.1	100	55-135
95-63-6	PS 1,2,4-Trimethylbenzene	50.0	0.00 U	48.9	98	53-132
135-98-8	PS sec-Butylbenzene	50.0	0.00 U	48.7	97	50-138
99-87-6	PS 4-Isopropyltoluene	50.0	0.00 U	48.9	98	49-138
541-73-1	PS 1,3-Dichlorobenzene	50.0	0.00 U	46.6	93	56-126
106-46-7	PS 1,4-Dichlorobenzene	50.0	0.00 U	46.7	93	55-125
104-51-8	PS n-Butylbenzene	50.0	0.00 U	45.9	92	43-142
96-12-8	PS 1,2-Dibromo-3-chloropropane	50.0	0.00 U	45.5	91	62-141
87-68-3	PS Hexachlorobutadiene	50.0	0.00 U	49.7	99	40-147
91-20-3	PS Naphthalene	50.0	0.00 U	48.9	98	62-134
87-61-6	PS 1,2,3-Trichlorobenzene	50.0	0.00 U	47.3	95	52-135

Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Post Spike

Client ID: CAWA-18-67PS

Matrix: W

Lab Sample ID 1203984437

Instrument: VOA9.I

Analysis Date: 03/08/2018 02:43

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No		Parmname	Amount Added ug/L	Sample Conc. ug/L		Spike Conc. ug/L	Recovery %	Acceptance Limits
120-82-1	PS	1,2,4-Trichlorobenzene	50.0	0.00	U	46.5	93	50-133
630-20-6	PS	1,1,1,2-Tetrachloroethane	50.0	0.00	U	52.4	105	71-133
95-50-1	PS	1,2-Dichlorobenzene	50.0	0.00	U	47.7	95	60-125
71-36-3	PS	n-Butyl alcohol	5000	0.00	U	4640	93	60-140

Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Post Spike Duplicate

Client ID: CAWA-18-67PSD

Matrix: W

Lab Sample ID 1203984439

Instrument: VOA9.I

Analysis Date: 03/08/2018 03:04

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits	RPD %	Acceptance Limits
179601-23-1	PSD m,p-Xylenes	100	0.00 U	108	108	59-132	10	0-20
75-05-8	PSD Acetonitrile	1250	0.00 U	1150	92	56-131	9	0-20
67-64-1	PSD Acetone	250	0.00 U	158	63	25-155	9	0-20
74-88-4	PSD Iodomethane	250	0.00 U	250	100	66-133	9	0-20
75-15-0	PSD Carbon disulfide	250	0.00 U	242	97	61-141	7	0-20
108-05-4	PSD Vinyl acetate	250	0.00 U	230	92	48-133	10	0-20
78-93-3	PSD 2-Butanone	250	0.00 U	202	81	25-143	11	0-20
108-10-1	PSD 4-Methyl-2-pentanone	250	0.00 U	246	98	61-127	13	0-20
591-78-6	PSD 2-Hexanone	250	0.00 U	217	87	33-138	11	0-20
75-71-8	PSD Dichlorodifluoromethane	50.0	0.00 U	54.3	109	33-164	9	0-20
74-87-3	PSD Chloromethane	50.0	0.00 U	65.2	130	53-139	9	0-20
75-01-4	PSD Vinyl chloride	50.0	0.00 U	65.4	131	58-140	9	0-20
74-83-9	PSD Bromomethane	50.0	0.00 U	54.6	109	59-146	9	0-20
75-00-3	PSD Chloroethane	50.0	0.00 U	54.6	109	65-129	8	0-20
75-69-4	PSD Trichlorofluoromethane	50.0	0.00 U	53.4	107	65-141	9	0-20
60-29-7	PSD Ethyl ether	50.0	0.00 U	56.0	112	69-127	11	0-20
75-35-4	PSD 1,1-Dichloroethylene	50.0	0.00 U	49.8	100	59-130	7	0-20
75-09-2	PSD Methylene chloride	50.0	0.00 U	49.6	99	62-123	10	0-20
1634-04-4	PSD tert-Butyl methyl ether	50.0	0.00 U	53.4	107	69-132	11	0-20
156-60-5	PSD trans-1,2-Dichloroethylene	50.0	0.00 U	51.2	102	65-127	9	0-20
75-34-3	PSD 1,1-Dichloroethane	50.0	0.00 U	52.7	105	67-127	10	0-20
156-59-2	PSD cis-1,2-Dichloroethylene	50.0	0.00 U	51.9	104	69-127	9	0-20

Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Post Spike Duplicate

Client ID: CAWA-18-67PSD

Matrix: W

Lab Sample ID 1203984439

Instrument: VOA9.I

Analysis Date: 03/08/2018 03:04

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits	RPD %	Acceptance Limits
594-20-7	PSD 2,2-Dichloropropane	50.0	0.00	U 48.9	98	66-137	9	0-20
74-97-5	PSD Bromochloromethane	50.0	0.00	U 54.1	108	71-130	12	0-20
67-66-3	PSD Chloroform	50.0	0.00	U 51.8	104	71-129	10	0-20
71-55-6	PSD 1,1,1-Trichloroethane	50.0	0.00	U 52.0	104	69-139	9	0-20
563-58-6	PSD 1,1-Dichloropropene	50.0	0.00	U 48.6	97	67-130	10	0-20
56-23-5	PSD Carbon tetrachloride	50.0	0.00	U 53.4	107	66-143	8	0-20
107-06-2	PSD 1,2-Dichloroethane	50.0	0.00	U 51.5	103	69-130	10	0-20
71-43-2	PSD Benzene	50.0	0.00	U 50.9	102	66-125	10	0-20
79-01-6	PSD Trichloroethylene	50.0	0.00	U 51.5	103	65-131	9	0-20
78-87-5	PSD 1,2-Dichloropropane	50.0	0.00	U 52.4	105	67-127	9	0-20
74-95-3	PSD Dibromomethane	50.0	0.00	U 53.2	106	72-129	11	0-20
75-27-4	PSD Bromodichloromethane	50.0	0.00	U 57.6	115	70-138	10	0-20
10061-01-5	PSD cis-1,3-Dichloropropylene	50.0	0.00	U 55.9	112	70-134	10	0-20
108-88-3	PSD Toluene	50.0	0.00	U 51.2	102	60-126	11	0-20
10061-02-6	PSD trans-1,3-Dichloropropylene	50.0	0.00	U 57.6	115	69-135	11	0-20
79-00-5	PSD 1,1,2-Trichloroethane	50.0	0.00	U 54.6	109	66-125	12	0-20
142-28-9	PSD 1,3-Dichloropropane	50.0	0.00	U 52.3	105	67-124	10	0-20
127-18-4	PSD Tetrachloroethylene	50.0	0.00	U 51.4	103	60-130	10	0-20
124-48-1	PSD Dibromochloromethane	50.0	0.00	U 56.0	112	68-143	13	0-20
106-93-4	PSD 1,2-Dibromoethane	50.0	0.00	U 56.1	112	71-127	10	0-20
108-90-7	PSD Chlorobenzene	50.0	0.00	U 52.8	106	64-124	11	0-20
100-41-4	PSD Ethylbenzene	50.0	0.00	U 51.2	102	61-130	9	0-20

Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Post Spike Duplicate

Client ID: CAWA-18-67PSD

Matrix: W

Lab Sample ID 1203984439

Instrument: VOA9.I

Analysis Date: 03/08/2018 03:04

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits	RPD %	Acceptance Limits
95-47-6	PSD o-Xylene	50.0	0.00	U 53.6	107	62-131	11	0-20
100-42-5	PSD Styrene	50.0	0.00	U 55.9	112	59-135	10	0-20
75-25-2	PSD Bromoform	50.0	0.00	U 66.3	133	64-138	11	0-20
98-82-8	PSD Isopropylbenzene	50.0	0.00	U 52.9	106	55-133	8	0-20
79-34-5	PSD 1,1,2,2-Tetrachloroethane	50.0	0.00	U 55.0	110	62-129	9	0-20
96-18-4	PSD 1,2,3-Trichloropropane	50.0	0.00	U 54.3	109	70-124	8	0-20
108-86-1	PSD Bromobenzene	50.0	0.00	U 54.0	108	62-124	9	0-20
103-65-1	PSD n-Propylbenzene	50.0	0.00	U 50.7	101	50-133	7	0-20
108-67-8	PSD 1,3,5-Trimethylbenzene	50.0	0.00	U 53.4	107	53-135	8	0-20
95-49-8	PSD 2-Chlorotoluene	50.0	0.00	U 53.8	108	56-128	8	0-20
106-43-4	PSD 4-Chlorotoluene	50.0	0.00	U 51.4	103	53-130	9	0-20
98-06-6	PSD tert-Butylbenzene	50.0	0.00	U 55.5	111	55-135	10	0-20
95-63-6	PSD 1,2,4-Trimethylbenzene	50.0	0.00	U 52.9	106	53-132	8	0-20
135-98-8	PSD sec-Butylbenzene	50.0	0.00	U 52.8	106	50-138	8	0-20
99-87-6	PSD 4-Isopropyltoluene	50.0	0.00	U 53.2	106	49-138	9	0-20
541-73-1	PSD 1,3-Dichlorobenzene	50.0	0.00	U 51.4	103	56-126	10	0-20
106-46-7	PSD 1,4-Dichlorobenzene	50.0	0.00	U 51.3	103	55-125	9	0-20
104-51-8	PSD n-Butylbenzene	50.0	0.00	U 50.0	100	43-142	9	0-20
96-12-8	PSD 1,2-Dibromo-3-chloropropane	50.0	0.00	U 52.2	104	62-141	14	0-20
87-68-3	PSD Hexachlorobutadiene	50.0	0.00	U 56.4	113	40-147	13	0-20
91-20-3	PSD Naphthalene	50.0	0.00	U 57.8	116	62-134	17	0-20
87-61-6	PSD 1,2,3-Trichlorobenzene	50.0	0.00	U 56.5	113	52-135	18	0-20

Volatile

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Quality Control Summary
Spike Recovery Report

SDG Number: 2018-1811

Sample Type: Post Spike Duplicate

Client ID: CAWA-18-67PSD

Matrix: W

Lab Sample ID 1203984439

Instrument: VOA9.I

Analysis Date: 03/08/2018 03:04

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No			Parmname		Amount Added ug/L	Sample Conc. ug/L	U	Spike Conc. ug/L	Recovery %	Acceptance Limits	Acceptance RPD %	Acceptance Limits
120-82-1	PSD	1,2,4-Trichlorobenzene			50.0	0.00	U	54.2	108	50-133	15	0-20
630-20-6	PSD	1,1,1,2-Tetrachloroethane			50.0	0.00	U	59.6	119	71-133	13	0-20
95-50-1	PSD	1,2-Dichlorobenzene			50.0	0.00	U	52.7	105	60-125	10	0-20
71-36-3	PSD	n-Butyl alcohol			5000	0.00	U	5180	104	60-140	11	0-20

Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Post Spike

Client ID: CAWA-18-67PS

Matrix: W

Lab Sample ID 1203984438

Instrument: VOA9.I

Analysis Date: 03/08/2018 03:33

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
107-02-8	PS Acrolein	250	0.00 U	232	93	49-141
76-13-1	PS Trichlorotrifluoroethane	250	0.00 U	244	98	57-149
107-05-1	PS Allyl chloride	250	0.00 U	233	93	54-128
107-13-1	PS Acrylonitrile	250	0.00 U	244	98	59-129
107-12-0	PS Propionitrile	250	0.00 U	233	93	58-131
126-98-7	PS Methacrylonitrile	250	0.00 U	238	95	59-134
80-62-6	PS Methyl methacrylate	250	0.00 U	251	101	62-135
97-63-2	PS Ethyl methacrylate	250	0.00 U	240	96	60-136
78-83-1	PS Isobutyl alcohol	2500	0.00 U	2360	94	60-143
126-99-8	PS 2-Chloro-1,3-butadiene	50.0	0.00 U	50.3	101	63-146

Volatile
Quality Control Summary
Spike Recovery Report

Page 2 of 2

SDG Number: 2018-1811

Sample Type: Post Spike Duplicate

Client ID: CAWA-18-67PSD

Matrix: W

Lab Sample ID 1203984440

Instrument: VOA9.I

Analysis Date: 03/08/2018 04:02

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	U	Spike Conc. ug/L	Recovery %	Acceptance Limits	RPD %	Acceptance Limits
107-02-8	PSD Acrolein	250	0.00	U	249	100	49-141	7	0-20
76-13-1	PSD Trichlorotrifluoroethane	250	0.00	U	252	101	57-149	3	0-20
107-05-1	PSD Allyl chloride	250	0.00	U	238	95	54-128	2	0-20
107-13-1	PSD Acrylonitrile	250	0.00	U	257	103	59-129	5	0-20
107-12-0	PSD Propionitrile	250	0.00	U	250	100	58-131	7	0-20
126-98-7	PSD Methacrylonitrile	250	0.00	U	251	100	59-134	5	0-20
80-62-6	PSD Methyl methacrylate	250	0.00	U	266	106	62-135	6	0-20
97-63-2	PSD Ethyl methacrylate	250	0.00	U	250	100	60-136	4	0-20
78-83-1	PSD Isobutyl alcohol	2500	0.00	U	2520	101	60-143	7	0-20
126-99-8	PSD 2-Chloro-1,3-butadiene	50.0	0.00	U	51.6	103	63-146	3	0-20

Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203985581

Instrument: VOA9.I

Analysis Date: 03/07/2018 23:45

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
179601-23-1	LCS m,p-Xylenes	100	0.0	80.7	81	71-127
75-05-8	LCS Acetonitrile	1250	0.0	979	78	61-125
67-64-1	LCS Acetone	250	0.0	158	63	48-157
74-88-4	LCS Iodomethane	250	0.0	197	79	72-128
75-15-0	LCS Carbon disulfide	250	0.0	192	77	69-138
108-05-4	LCS Vinyl acetate	250	0.0	227	91	67-125
78-93-3	LCS 2-Butanone	250	0.0	177	71	55-138
108-10-1	LCS 4-Methyl-2-pentanone	250	0.0	201	81	66-124
591-78-6	LCS 2-Hexanone	250	0.0	179	72	56-140
75-71-8	LCS Dichlorodifluoromethane	50.0	0.0	46.3	93	40-160
74-87-3	LCS Chloromethane	50.0	0.0	58.4	117	58-135
75-01-4	LCS Vinyl chloride	50.0	0.0	58.8	118	65-137
74-83-9	LCS Bromomethane	50.0	0.0	48.8	98	63-137
75-00-3	LCS Chloroethane	50.0	0.0	47.3	95	69-129
75-69-4	LCS Trichlorofluoromethane	50.0	0.0	47.4	95	69-138
60-29-7	LCS Ethyl ether	50.0	0.0	50.9	102	72-125
75-35-4	LCS 1,1-Dichloroethylene	50.0	0.0	38.8	78	66-126
75-09-2	LCS Methylene chloride	50.0	0.0	39.6	79	68-119
1634-04-4	LCS tert-Butyl methyl ether	50.0	0.0	42.4	85	76-128
156-60-5	LCS trans-1,2-Dichloroethylene	50.0	0.0	40.6	81	71-124
75-34-3	LCS 1,1-Dichloroethane	50.0	0.0	41.7	83	73-123
156-59-2	LCS cis-1,2-Dichloroethylene	50.0	0.0	41.8	84	75-123

Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203985581

Instrument: VOA9.I

Analysis Date: 03/07/2018 23:45

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
594-20-7	LCS 2,2-Dichloropropane	50.0	0.0	39.3	79	72-138
74-97-5	LCS Bromochloromethane	50.0	0.0	43.4	87	76-125
67-66-3	LCS Chloroform	50.0	0.0	41.5	83	76-123
71-55-6	LCS 1,1,1-Trichloroethane	50.0	0.0	40.1	80	74-136
563-58-6	LCS 1,1-Dichloropropene	50.0	0.0	37.0	74	72-129
56-23-5	LCS Carbon tetrachloride	50.0	0.0	40.3	81	72-140
107-06-2	LCS 1,2-Dichloroethane	50.0	0.0	41.8	84	74-122
71-43-2	LCS Benzene	50.0	0.0	40.1	80	72-121
79-01-6	LCS Trichloroethylene	50.0	0.0	39.5	79	74-125
78-87-5	LCS 1,2-Dichloropropane	50.0	0.0	42.0	84	73-121
74-95-3	LCS Dibromomethane	50.0	0.0	42.2	84	78-123
75-27-4	LCS Bromodichloromethane	50.0	0.0	45.2	90	77-131
10061-01-5	LCS cis-1,3-Dichloropropylene	50.0	0.0	44.3	89	78-131
108-88-3	LCS Toluene	50.0	0.0	39.9	80	71-121
10061-02-6	LCS trans-1,3-Dichloropropylene	50.0	0.0	46.2	92	78-131
79-00-5	LCS 1,1,2-Trichloroethane	50.0	0.0	43.5	87	74-118
142-28-9	LCS 1,3-Dichloropropane	50.0	0.0	42.1	84	74-118
127-18-4	LCS Tetrachloroethylene	50.0	0.0	36.8	74	69-129
124-48-1	LCS Dibromochloromethane	50.0	0.0	43.1	86	76-137
106-93-4	LCS 1,2-Dibromoethane	50.0	0.0	44.6	89	78-122
108-90-7	LCS Chlorobenzene	50.0	0.0	40.6	81	74-120
100-41-4	LCS Ethylbenzene	50.0	0.0	38.8	78	73-125

Volatile
Quality Control Summary
Spike Recovery Report

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203985581

Instrument: VOA9.I

Analysis Date: 03/07/2018 23:45

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
95-47-6	LCS o-Xylene	50.0	0.0	40.6	81	74-126
100-42-5	LCS Styrene	50.0	0.0	43.2	86	72-130
75-25-2	LCS Bromoform	50.0	0.0	52.3	105	72-136
98-82-8	LCS Isopropylbenzene	50.0	0.0	39.3	79	70-130
79-34-5	LCS 1,1,2,2-Tetrachloroethane	50.0	0.0	44.9	90	70-126
96-18-4	LCS 1,2,3-Trichloropropane	50.0	0.0	45.1	90	74-122
108-86-1	LCS Bromobenzene	50.0	0.0	42.7	85	74-120
103-65-1	LCS n-Propylbenzene	50.0	0.0	37.4	75	67-128
108-67-8	LCS 1,3,5-Trimethylbenzene	50.0	0.0	39.3	79	70-129
95-49-8	LCS 2-Chlorotoluene	50.0	0.0	40.5	81	71-124
106-43-4	LCS 4-Chlorotoluene	50.0	0.0	39.6	79	69-125
98-06-6	LCS tert-Butylbenzene	50.0	0.0	38.0	76	72-130
95-63-6	LCS 1,2,4-Trimethylbenzene	50.0	0.0	39.5	79	70-126
135-98-8	LCS sec-Butylbenzene	50.0	0.0	36.3	73	70-131
99-87-6	LCS 4-Isopropyltoluene	50.0	0.0	36.6	73	71-131
541-73-1	LCS 1,3-Dichlorobenzene	50.0	0.0	39.2	78	72-121
106-46-7	LCS 1,4-Dichlorobenzene	50.0	0.0	39.8	80	71-120
104-51-8	LCS n-Butylbenzene	50.0	0.0	33.8	68	68-134
96-12-8	LCS 1,2-Dibromo-3-chloropropane	50.0	0.0	42.5	85	68-141
87-68-3	LCS Hexachlorobutadiene	50.0	0.0	33.2	66 *	72-136
91-20-3	LCS Naphthalene	50.0	0.0	46.3	93	72-132
87-61-6	LCS 1,2,3-Trichlorobenzene	50.0	0.0	42.5	85	70-130

Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203985581

Instrument: VOA9.I

Analysis Date: 03/07/2018 23:45

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
120-82-1	LCS 1,2,4-Trichlorobenzene	50.0	0.0	39.4	79	71-129
630-20-6	LCS 1,1,1,2-Tetrachloroethane	50.0	0.0	45.6	91	79-127
95-50-1	LCS 1,2-Dichlorobenzene	50.0	0.0	40.9	82	74-120
71-36-3	LCS n-Butyl alcohol	5000	0.0	4220	84	63-138

Volatile
Quality Control Summary
Spike Recovery Report

Page 1 of 1

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1744716

Matrix: WATER

Lab Sample ID 1203985582

Instrument: VOA9.I

Analysis Date: 03/08/2018 00:41

Dilution: 1

Analyst: RXY1

Purge Vol: 5 mL

Batch ID: 1744716

CAS No		Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
107-02-8	LCS	Acrolein	250	0.0	236	94	60-140
76-13-1	LCS	Trichlorotrifluoroethane	250	0.0	211	84	61-148
107-05-1	LCS	Allyl chloride	250	0.0	214	85	59-125
107-13-1	LCS	Acrylonitrile	250	0.0	234	93	65-122
107-12-0	LCS	Propionitrile	250	0.0	227	91	64-124
126-98-7	LCS	Methacrylonitrile	250	0.0	226	90	64-126
80-62-6	LCS	Methyl methacrylate	250	0.0	232	93	69-127
97-63-2	LCS	Ethyl methacrylate	250	0.0	225	90	66-130
78-83-1	LCS	Isobutyl alcohol	2500	0.0	2330	93	65-135
126-99-8	LCS	2-Chloro-1,3-butadiene	50.0	0.0	44.7	89	66-147

Method Blank Summary

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SDG Number: 2018-1811

Client: ARSL004

Matrix: WATER

Client ID: MB for batch 1744716

Instrument ID: VOA9.I

Data File: 030618V9\9I230B1.D

Lab Sample ID: 1203984431

Prep Date: 03/07/2018 01:40

Analyzed: 03/07/18 01:40

Column: DB-624

This method blank applies to the following samples and quality control samples:

Client Sample ID	Lab Sample ID	File ID	Date Analyzed	Time Analyzed
01 LCS for batch 1744716	1203984433	030618V9\9I227L1.D	03/07/18	0014
02 LCS for batch 1744716	1203984434	030618V9\9I229L1.D	03/07/18	0111
03 CAPA-18-6	444566004	030618V9\9I231.D	03/07/18	0209
04 CAWA-18-111	444566010	030618V9\9I232.D	03/07/18	0238
05 CAWA-18-112	444566014	030618V9\9I233.D	03/07/18	0307
06 CAPA-18-2	444566002	030618V9\9I234.D	03/07/18	0336
07 CAPA-18-12	444566006	030618V9\9I235.D	03/07/18	0404
08 CAWA-18-54	444566009	030618V9\9I236.D	03/07/18	0433
09 CAWA-18-56	444566012	030618V9\9I237.D	03/07/18	0501

Method Blank Summary

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SDG Number:	2018-1811	Client:	ARSL004	Matrix:	WATER
Client ID:	MB for batch 1744716	Instrument ID:	VOA9.I	Data File:	030718V9\9I332B.D
Lab Sample ID:	1203985580	Prep Date:	03/08/2018 01:09	Analyzed:	03/08/18 01:09
Column:	DB-624				

This method blank applies to the following samples and quality control samples:

Client Sample ID	Lab Sample ID	File ID	Date Analyzed	Time Analyzed
11 LCS for batch 1744716	1203985581	030718V9\9I329L.D	03/07/18	2345
12 LCS for batch 1744716	1203985582	030718V9\9I331L.D	03/08/18	0041
13 CAWA-18-67PS	1203984437	030718V9\9I335.D	03/08/18	0243
14 CAWA-18-67PSD	1203984439	030718V9\9I336.D	03/08/18	0304
15 CAWA-18-67PS	1203984438	030718V9\9I337.D	03/08/18	0333
16 CAWA-18-67PSD	1203984440	030718V9\9I338.D	03/08/18	0402

Quality Control Data

Volatile
Certificate of Analysis
Sample Summary

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SDG Number: 2018-1811

Lab Sample ID: 1203984431

Client Sample: QC for batch 1744716

Client ID: MB for batch 1744716

Batch ID: 1744716

Run Date: 03/07/2018 01:40

Prep Date: 03/07/2018 01:40

Data File: 030618V9\9I230B1.D

Client: ARSL004

Method: SW-846:8260B

Inst: VOA9.I

Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC

SOP Ref: GL-OA-E-038

Dilution: 1

Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

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SDG Number: 2018-1811
Lab Sample ID: 1203984431
Client Sample: QC for batch 1744716
Client ID: MB for batch 1744716
Batch ID: 1744716
Run Date: 03/07/2018 01:40
Prep Date: 03/07/2018 01:40
Data File: 030618V9\91230B1.D

Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811
Lab Sample ID: 1203984431
Client Sample: QC for batch 1744716
Client ID: MB for batch 1744716
Batch ID: 1744716
Run Date: 03/07/2018 01:40
Prep Date: 03/07/2018 01:40
Data File: 030618V9\9I230B1.D

Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	49.0	50.0	ug/L 98	(71%-134%)
Bromofluorobenzene	49.4	50.0	ug/L 99	(70%-131%)
Toluene-d8	50.7	50.0	ug/L 101	(74%-124%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
	unknown siloxane	18.968	39.6	ug/L	0	J

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 1203984433

Client Sample: QC for batch 1744716

Client ID: LCS for batch 1744716

Batch ID: 1744716

Run Date: 03/07/2018 00:14

Prep Date: 03/07/2018 00:14

Data File: 030618V9\9I227L1.D

Client: ARSL004

Method: SW-846:8260B

Inst: VOA9.I

Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC

SOP Ref: GL-OA-E-038

Dilution: 1

Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane		46.5	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane		40.0	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane		48.7	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane		46.0	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane		41.5	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene		38.8	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene		37.5	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene		47.3	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane		49.3	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene		44.2	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene		41.8	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane		47.9	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane		47.3	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene		42.9	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane		42.8	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane		41.7	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene		42.2	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene		41.2	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane		44.1	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene		41.5	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane		38.5	ug/L	0.300	1.00
78-93-3	2-Butanone		215	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene		42.0	ug/L	0.300	1.00
591-78-6	2-Hexanone		212	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene		40.5	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene		41.0	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone		233	ug/L	1.50	5.00
67-64-1	Acetone		198	ug/L	1.50	10.0
75-05-8	Acetonitrile		1120	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene		40.1	ug/L	0.300	1.00
108-86-1	Bromobenzene		43.3	ug/L	0.300	1.00
74-97-5	Bromochloromethane		44.9	ug/L	0.300	1.00
75-27-4	Bromodichloromethane		45.6	ug/L	0.300	1.00
75-25-2	Bromoform		55.7	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

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SDG Number: 2018-1811
Lab Sample ID: 1203984433
Client Sample: QC for batch 1744716
Client ID: LCS for batch 1744716
Batch ID: 1744716
Run Date: 03/07/2018 00:14
Prep Date: 03/07/2018 00:14
Data File: 030618V9\9I227L1.D

Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane		40.9	ug/L	0.300	1.00
75-15-0	Carbon disulfide		193	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride		40.4	ug/L	0.300	1.00
108-90-7	Chlorobenzene		41.7	ug/L	0.300	1.00
75-00-3	Chloroethane		41.9	ug/L	0.300	1.00
67-66-3	Chloroform		41.5	ug/L	0.300	1.00
74-87-3	Chloromethane		46.8	ug/L	0.300	1.00
124-48-1	Dibromochloromethane		45.6	ug/L	0.300	1.00
74-95-3	Dibromomethane		44.5	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane		42.2	ug/L	0.300	1.00
60-29-7	Ethyl ether		48.4	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene		40.3	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene		40.7	ug/L	0.300	1.00
74-88-4	Iodomethane		201	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene		41.5	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride		39.9	ug/L	1.00	10.0
91-20-3	Naphthalene		52.0	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene		44.8	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene		39.4	ug/L	0.300	1.00
108-88-3	Toluene		40.8	ug/L	0.300	1.00
79-01-6	Trichloroethylene		40.1	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane		41.5	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate		233	ug/L	1.50	5.00
75-01-4	Vinyl chloride		47.9	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene		41.3	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene		44.9	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes		83.9	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol		5020	ug/L	15.0	50.0
104-51-8	n-Butylbenzene		38.4	ug/L	0.300	1.00
103-65-1	n-Propylbenzene		39.7	ug/L	0.300	1.00
95-47-6	o-Xylene		42.1	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene		40.7	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811
Lab Sample ID: 1203984433
Client Sample: QC for batch 1744716
Client ID: LCS for batch 1744716
Batch ID: 1744716
Run Date: 03/07/2018 00:14
Prep Date: 03/07/2018 00:14
Data File: 030618V9\9I227L1.D

Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether		45.6	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene		41.8	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene		40.1	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene		48.2	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	49.6	50.0	ug/L	99	(71%-134%)
Bromofluorobenzene	49.4	50.0	ug/L	99	(70%-131%)
Toluene-d8	51.6	50.0	ug/L	103	(74%-124%)

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 1203984434

Client Sample: QC for batch 1744716

Client ID: LCS for batch 1744716

Batch ID: 1744716

Run Date: 03/07/2018 01:11

Prep Date: 03/07/2018 01:11

Data File: 030618V9\9I229L1.D

Client: ARSL004

Method: SW-846:8260B

Inst: VOA9.I

Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC

SOP Ref: GL-OA-E-038

Dilution: 1

Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene		47.4	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein		251	ug/L	1.50	5.00
107-13-1	Acrylonitrile		259	ug/L	1.50	5.00
107-05-1	Allyl chloride		222	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811
Lab Sample ID: 1203984434
Client Sample: QC for batch 1744716
Client ID: LCS for batch 1744716
Batch ID: 1744716
Run Date: 03/07/2018 01:11
Prep Date: 03/07/2018 01:11
Data File: 030618V9\9I229L1.D

Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate		242	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol		2630	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile		248	ug/L	1.50	5.00
80-62-6	Methyl methacrylate		251	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile		255	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane		227	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

SDG Number:	2018-1811	Matrix:	WATER
Lab Sample ID:	1203984434		
Client Sample:	QC for batch 1744716	Client:	ARSL004
Client ID:	LCS for batch 1744716	Method:	SW-846:8260B
Batch ID:	1744716	Inst:	VOA9.I
Run Date:	03/07/2018 01:11	Analyst:	RXY1
Prep Date:	03/07/2018 01:11		
Data File:	030618V9\9I229L1.D	Column:	DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	50.2	50.0	ug/L	100	(71%-134%)
Bromofluorobenzene	51.4	50.0	ug/L	103	(70%-131%)
Toluene-d8	53.0	50.0	ug/L	106	(74%-124%)

Volatile
Certificate of Analysis
Sample Summary

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SDG Number: 2018-1811
Lab Sample ID: 1203984437
Client Sample: QC for batch 1744716
Client ID: CAWA-18-67PS
Batch ID: 1744716
Run Date: 03/08/2018 02:43
Prep Date: 03/08/2018 02:43
Data File: 030718V9\9I335.D

Date Collected: 02/22/2018 13:43
Date Received: 02/24/2018 09:25
Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: W

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane		52.4	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane		47.6	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane		50.3	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane		48.3	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane		47.6	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene		46.4	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene		44.1	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene		47.3	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane		50.0	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene		46.5	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene		48.9	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane		45.5	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane		50.6	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene		47.7	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane		46.6	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane		47.6	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene		49.4	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene		46.6	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane		47.4	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene		46.7	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane		44.6	ug/L	0.300	1.00
78-93-3	2-Butanone		181	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene		49.4	ug/L	0.300	1.00
591-78-6	2-Hexanone		194	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene		47.2	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene		48.9	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone		216	ug/L	1.50	5.00
67-64-1	Acetone		144	ug/L	1.50	10.0
75-05-8	Acetonitrile		1050	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene		46.3	ug/L	0.300	1.00
108-86-1	Bromobenzene		49.3	ug/L	0.300	1.00
74-97-5	Bromochloromethane		48.2	ug/L	0.300	1.00
75-27-4	Bromodichloromethane		52.1	ug/L	0.300	1.00
75-25-2	Bromoform		59.4	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

SDG Number: 2018-1811	Date Collected: 02/22/2018 13:43	Matrix: W
Lab Sample ID: 1203984437	Date Received: 02/24/2018 09:25	
Client Sample: QC for batch 1744716	Client: ARSL004	Project: QC
Client ID: CAWA-18-67PS	Method: SW-846:8260B	SOP Ref: GL-OA-E-038
Batch ID: 1744716	Inst: VOA9.I	Dilution: 1
Run Date: 03/08/2018 02:43	Analyst: RXY1	Purge Vol: 5 mL
Prep Date: 03/08/2018 02:43		
Data File: 030718V9\9I335.D	Column: DB-624	

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane		50.1	ug/L	0.300	1.00
75-15-0	Carbon disulfide		226	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride		49.1	ug/L	0.300	1.00
108-90-7	Chlorobenzene		47.3	ug/L	0.300	1.00
75-00-3	Chloroethane		50.5	ug/L	0.300	1.00
67-66-3	Chloroform		46.6	ug/L	0.300	1.00
74-87-3	Chloromethane		59.6	ug/L	0.300	1.00
124-48-1	Dibromochloromethane		49.3	ug/L	0.300	1.00
74-95-3	Dibromomethane		47.9	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane		49.5	ug/L	0.300	1.00
60-29-7	Ethyl ether		50.3	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene		46.7	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene		49.7	ug/L	0.300	1.00
74-88-4	Iodomethane		228	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene		48.6	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride		45.0	ug/L	1.00	10.0
91-20-3	Naphthalene		48.9	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene		50.5	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene		46.5	ug/L	0.300	1.00
108-88-3	Toluene		45.9	ug/L	0.300	1.00
79-01-6	Trichloroethylene		46.9	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane		49.0	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate		209	ug/L	1.50	5.00
75-01-4	Vinyl chloride		59.8	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene		47.2	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene		50.4	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes		97.6	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol		4640	ug/L	15.0	50.0
104-51-8	n-Butylbenzene		45.9	ug/L	0.300	1.00
103-65-1	n-Propylbenzene		47.4	ug/L	0.300	1.00
95-47-6	o-Xylene		48.0	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene		48.7	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

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SDG Number:	2018-1811	Date Collected:	02/22/2018 13:43	Matrix:	W
Lab Sample ID:	1203984437	Date Received:	02/24/2018 09:25		
Client Sample:	QC for batch 1744716	Client:	ARSL004	Project:	QC
Client ID:	CAWA-18-67PS	Method:	SW-846:8260B	SOP Ref:	GL-OA-E-038
Batch ID:	1744716	Inst:	VOA9.I	Dilution:	1
Run Date:	03/08/2018 02:43	Analyst:	RXY1	Purge Vol:	5 mL
Prep Date:	03/08/2018 02:43				
Data File:	030718V9\9I335.D	Column:	DB-624		

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether		47.9	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene		50.1	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene		46.9	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene		51.4	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	49.9	50.0	ug/L	100	(71%-134%)
Bromofluorobenzene	51.3	50.0	ug/L	103	(70%-131%)
Toluene-d8	51.4	50.0	ug/L	103	(74%-124%)

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811
Lab Sample ID: 1203984438
Client Sample: QC for batch 1744716
Client ID: CAWA-18-67PS
Batch ID: 1744716
Run Date: 03/08/2018 03:33
Prep Date: 03/08/2018 03:33
Data File: 030718V9\9I337.D

Date Collected: 02/22/2018 13:43
Date Received: 02/24/2018 09:25
Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: W

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene		50.3	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein		232	ug/L	1.50	5.00
107-13-1	Acrylonitrile		244	ug/L	1.50	5.00
107-05-1	Allyl chloride		233	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

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SDG Number: 2018-1811
Lab Sample ID: 1203984438
Client Sample: QC for batch 1744716
Client ID: CAWA-18-67PS
Batch ID: 1744716
Run Date: 03/08/2018 03:33
Prep Date: 03/08/2018 03:33
Data File: 030718V9\9I337.D

Date Collected: 02/22/2018 13:43
Date Received: 02/24/2018 09:25
Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: W

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate		240	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol		2360	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile		238	ug/L	1.50	5.00
80-62-6	Methyl methacrylate		251	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile		233	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane		244	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

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SDG Number:	2018-1811	Date Collected:	02/22/2018 13:43	Matrix:	W
Lab Sample ID:	1203984438	Date Received:	02/24/2018 09:25		
Client Sample:	QC for batch 1744716	Client:	ARSL004	Project:	QC
Client ID:	CAWA-18-67PS	Method:	SW-846:8260B	SOP Ref:	GL-OA-E-038
Batch ID:	1744716	Inst:	VOA9.I	Dilution:	1
Run Date:	03/08/2018 03:33	Analyst:	RXY1	Purge Vol:	5 mL
Prep Date:	03/08/2018 03:33				
Data File:	030718V9\9I337.D	Column:	DB-624		

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	48.7	50.0	ug/L	97	(71%-134%)
Bromofluorobenzene	48.7	50.0	ug/L	97	(70%-131%)
Toluene-d8	50.8	50.0	ug/L	102	(74%-124%)

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SDG Number: 2018-1811	Date Collected: 02/22/2018 13:43	Matrix: W
Lab Sample ID: 1203984439	Date Received: 02/24/2018 09:25	
Client Sample: QC for batch 1744716	Client: ARSL004	Project: QC
Client ID: CAWA-18-67PSD	Method: SW-846:8260B	SOP Ref: GL-OA-E-038
Batch ID: 1744716	Inst: VOA9.I	Dilution: 1
Run Date: 03/08/2018 03:04	Analyst: RXY1	Purge Vol: 5 mL
Prep Date: 03/08/2018 03:04		
Data File: 030718V9\9I336.D	Column: DB-624	

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane		59.6	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane		52.0	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane		55.0	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane		54.6	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane		52.7	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene		49.8	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene		48.6	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene		56.5	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane		54.3	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene		54.2	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene		52.9	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane		52.2	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane		56.1	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene		52.7	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane		51.5	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane		52.4	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene		53.4	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene		51.4	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane		52.3	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene		51.3	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane		48.9	ug/L	0.300	1.00
78-93-3	2-Butanone		202	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene		53.8	ug/L	0.300	1.00
591-78-6	2-Hexanone		217	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene		51.4	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene		53.2	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone		246	ug/L	1.50	5.00
67-64-1	Acetone		158	ug/L	1.50	10.0
75-05-8	Acetonitrile		1150	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene		50.9	ug/L	0.300	1.00
108-86-1	Bromobenzene		54.0	ug/L	0.300	1.00
74-97-5	Bromochloromethane		54.1	ug/L	0.300	1.00
75-27-4	Bromodichloromethane		57.6	ug/L	0.300	1.00
75-25-2	Bromoform		66.3	ug/L	0.300	1.00

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Sample Summary

SDG Number: 2018-1811	Date Collected: 02/22/2018 13:43	Matrix: W
Lab Sample ID: 1203984439	Date Received: 02/24/2018 09:25	
Client Sample: QC for batch 1744716	Client: ARSL004	Project: QC
Client ID: CAWA-18-67PSD	Method: SW-846:8260B	SOP Ref: GL-OA-E-038
Batch ID: 1744716	Inst: VOA9.I	Dilution: 1
Run Date: 03/08/2018 03:04	Analyst: RXY1	Purge Vol: 5 mL
Prep Date: 03/08/2018 03:04		
Data File: 030718V9\9I336.D	Column: DB-624	

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane		54.6	ug/L	0.300	1.00
75-15-0	Carbon disulfide		242	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride		53.4	ug/L	0.300	1.00
108-90-7	Chlorobenzene		52.8	ug/L	0.300	1.00
75-00-3	Chloroethane		54.6	ug/L	0.300	1.00
67-66-3	Chloroform		51.8	ug/L	0.300	1.00
74-87-3	Chloromethane		65.2	ug/L	0.300	1.00
124-48-1	Dibromochloromethane		56.0	ug/L	0.300	1.00
74-95-3	Dibromomethane		53.2	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane		54.3	ug/L	0.300	1.00
60-29-7	Ethyl ether		56.0	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene		51.2	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene		56.4	ug/L	0.300	1.00
74-88-4	Iodomethane		250	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene		52.9	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride		49.6	ug/L	1.00	10.0
91-20-3	Naphthalene		57.8	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene		55.9	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene		51.4	ug/L	0.300	1.00
108-88-3	Toluene		51.2	ug/L	0.300	1.00
79-01-6	Trichloroethylene		51.5	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane		53.4	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate		230	ug/L	1.50	5.00
75-01-4	Vinyl chloride		65.4	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene		51.9	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene		55.9	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes		108	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol		5180	ug/L	15.0	50.0
104-51-8	n-Butylbenzene		50.0	ug/L	0.300	1.00
103-65-1	n-Propylbenzene		50.7	ug/L	0.300	1.00
95-47-6	o-Xylene		53.6	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene		52.8	ug/L	0.300	1.00

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Sample Summary

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SDG Number:	2018-1811	Date Collected:	02/22/2018 13:43	Matrix:	W
Lab Sample ID:	1203984439	Date Received:	02/24/2018 09:25		
Client Sample:	QC for batch 1744716	Client:	ARSL004	Project:	QC
Client ID:	CAWA-18-67PSD	Method:	SW-846:8260B	SOP Ref:	GL-OA-E-038
Batch ID:	1744716	Inst:	VOA9.I	Dilution:	1
Run Date:	03/08/2018 03:04	Analyst:	RXY1	Purge Vol:	5 mL
Prep Date:	03/08/2018 03:04				
Data File:	030718V9\9I336.D	Column:	DB-624		

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether		53.4	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene		55.5	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene		51.2	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene		57.6	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	51.3	50.0	ug/L	103	(71%-134%)
Bromofluorobenzene	52.0	50.0	ug/L	104	(70%-131%)
Toluene-d8	53.3	50.0	ug/L	107	(74%-124%)

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Sample Summary

SDG Number: 2018-1811	Date Collected: 02/22/2018 13:43	Matrix: W
Lab Sample ID: 1203984440	Date Received: 02/24/2018 09:25	
Client Sample: QC for batch 1744716	Client: ARSL004	Project: QC
Client ID: CAWA-18-67PSD	Method: SW-846:8260B	SOP Ref: GL-OA-E-038
Batch ID: 1744716	Inst: VOA9.I	Dilution: 1
Run Date: 03/08/2018 04:02	Analyst: RXY1	Purge Vol: 5 mL
Prep Date: 03/08/2018 04:02		
Data File: 030718V9\9I338.D	Column: DB-624	

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene		51.6	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein		249	ug/L	1.50	5.00
107-13-1	Acrylonitrile		257	ug/L	1.50	5.00
107-05-1	Allyl chloride		238	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

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Sample Summary

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SDG Number: 2018-1811
Lab Sample ID: 1203984440
Client Sample: QC for batch 1744716
Client ID: CAWA-18-67PSD
Batch ID: 1744716
Run Date: 03/08/2018 04:02
Prep Date: 03/08/2018 04:02
Data File: 030718V9\9I338.D

Date Collected: 02/22/2018 13:43
Date Received: 02/24/2018 09:25
Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: W

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate		250	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol		2520	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile		251	ug/L	1.50	5.00
80-62-6	Methyl methacrylate		266	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile		250	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane		252	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

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SDG Number:	2018-1811	Date Collected:	02/22/2018 13:43	Matrix:	W
Lab Sample ID:	1203984440	Date Received:	02/24/2018 09:25		
Client Sample:	QC for batch 1744716	Client:	ARSL004	Project:	QC
Client ID:	CAWA-18-67PSD	Method:	SW-846:8260B	SOP Ref:	GL-OA-E-038
Batch ID:	1744716	Inst:	VOA9.I	Dilution:	1
Run Date:	03/08/2018 04:02	Analyst:	RXY1	Purge Vol:	5 mL
Prep Date:	03/08/2018 04:02				
Data File:	030718V9\9I338.D	Column:	DB-624		

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	49.1	50.0	ug/L	98	(71%-134%)
Bromofluorobenzene	50.5	50.0	ug/L	101	(70%-131%)
Toluene-d8	51.7	50.0	ug/L	103	(74%-124%)

Volatile
Certificate of Analysis
Sample Summary

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SDG Number: 2018-1811

Lab Sample ID: 1203985580

Client Sample: QC for batch 1744716

Client ID: MB for batch 1744716

Batch ID: 1744716

Run Date: 03/08/2018 01:09

Prep Date: 03/08/2018 01:09

Data File: 030718V9\9I332B.D

Client: ARSL004

Method: SW-846:8260B

Inst: VOA9.I

Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC

SOP Ref: GL-OA-E-038

Dilution: 1

Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

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SDG Number: 2018-1811
Lab Sample ID: 1203985580
Client Sample: QC for batch 1744716
Client ID: MB for batch 1744716
Batch ID: 1744716
Run Date: 03/08/2018 01:09
Prep Date: 03/08/2018 01:09
Data File: 030718V9\9I332B.D

Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

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SDG Number:	2018-1811	Matrix:	WATER
Lab Sample ID:	1203985580		
Client Sample:	QC for batch 1744716	Client:	ARSL004
Client ID:	MB for batch 1744716	Method:	SW-846:8260B
Batch ID:	1744716	Inst:	VOA9.I
Run Date:	03/08/2018 01:09	Analyst:	RXY1
Prep Date:	03/08/2018 01:09		
Data File:	030718V9\9I332B.D	Column:	DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	49.7	50.0	ug/L 99	(71%-134%)
Bromofluorobenzene	50.1	50.0	ug/L 100	(70%-131%)
Toluene-d8	51.4	50.0	ug/L 103	(74%-124%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
No Tentatively Identified Compounds Found				ug/L		

Volatile
Certificate of Analysis
Sample Summary

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SDG Number: 2018-1811

Lab Sample ID: 1203985581

Client Sample: QC for batch 1744716

Client ID: LCS for batch 1744716

Batch ID: 1744716

Run Date: 03/07/2018 23:45

Prep Date: 03/07/2018 23:45

Data File: 030718V9\9I329L.D

Client: ARSL004

Method: SW-846:8260B

Inst: VOA9.I

Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC

SOP Ref: GL-OA-E-038

Dilution: 1

Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane		45.6	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane		40.1	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane		44.9	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane		43.5	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane		41.7	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene		38.8	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene		37.0	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene		42.5	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane		45.1	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene		39.4	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene		39.5	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane		42.5	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane		44.6	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene		40.9	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane		41.8	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane		42.0	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene		39.3	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene		39.2	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane		42.1	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene		39.8	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane		39.3	ug/L	0.300	1.00
78-93-3	2-Butanone		177	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene	U	0.300	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene		40.5	ug/L	0.300	1.00
591-78-6	2-Hexanone		179	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene		39.6	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene		36.6	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone		201	ug/L	1.50	5.00
67-64-1	Acetone		158	ug/L	1.50	10.0
75-05-8	Acetonitrile		979	ug/L	8.00	25.0
107-02-8	Acrolein	U	1.50	ug/L	1.50	5.00
107-13-1	Acrylonitrile	U	1.50	ug/L	1.50	5.00
107-05-1	Allyl chloride	U	1.50	ug/L	1.50	5.00
71-43-2	Benzene		40.1	ug/L	0.300	1.00
108-86-1	Bromobenzene		42.7	ug/L	0.300	1.00
74-97-5	Bromochloromethane		43.4	ug/L	0.300	1.00
75-27-4	Bromodichloromethane		45.2	ug/L	0.300	1.00
75-25-2	Bromoform		52.3	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811
Lab Sample ID: 1203985581
Client Sample: QC for batch 1744716
Client ID: LCS for batch 1744716
Batch ID: 1744716
Run Date: 03/07/2018 23:45
Prep Date: 03/07/2018 23:45
Data File: 030718V9\9I329L.D

Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane		48.8	ug/L	0.300	1.00
75-15-0	Carbon disulfide		192	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride		40.3	ug/L	0.300	1.00
108-90-7	Chlorobenzene		40.6	ug/L	0.300	1.00
75-00-3	Chloroethane		47.3	ug/L	0.300	1.00
67-66-3	Chloroform		41.5	ug/L	0.300	1.00
74-87-3	Chloromethane		58.4	ug/L	0.300	1.00
124-48-1	Dibromochloromethane		43.1	ug/L	0.300	1.00
74-95-3	Dibromomethane		42.2	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane		46.3	ug/L	0.300	1.00
60-29-7	Ethyl ether		50.9	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate	U	1.50	ug/L	1.50	5.00
100-41-4	Ethylbenzene		38.8	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene		33.2	ug/L	0.300	1.00
74-88-4	Iodomethane		197	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol	U	15.0	ug/L	15.0	50.0
98-82-8	Isopropylbenzene		39.3	ug/L	0.300	1.00
126-98-7	Methacrylonitrile	U	1.50	ug/L	1.50	5.00
80-62-6	Methyl methacrylate	U	1.50	ug/L	1.50	5.00
75-09-2	Methylene chloride		39.6	ug/L	1.00	10.0
91-20-3	Naphthalene		46.3	ug/L	0.300	1.00
107-12-0	Propionitrile	U	1.50	ug/L	1.50	5.00
100-42-5	Styrene		43.2	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene		36.8	ug/L	0.300	1.00
108-88-3	Toluene		39.9	ug/L	0.300	1.00
79-01-6	Trichloroethylene		39.5	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane		47.4	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane	U	2.00	ug/L	2.00	5.00
108-05-4	Vinyl acetate		227	ug/L	1.50	5.00
75-01-4	Vinyl chloride		58.8	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene		41.8	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene		44.3	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes		80.7	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol		4220	ug/L	15.0	50.0
104-51-8	n-Butylbenzene		33.8	ug/L	0.300	1.00
103-65-1	n-Propylbenzene		37.4	ug/L	0.300	1.00
95-47-6	o-Xylene		40.6	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene		36.3	ug/L	0.300	1.00

**Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811
Lab Sample ID: 1203985581
Client Sample: QC for batch 1744716
Client ID: LCS for batch 1744716
Batch ID: 1744716
Run Date: 03/07/2018 23:45
Prep Date: 03/07/2018 23:45
Data File: 030718V9\9I329L.D

Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether		42.4	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene		38.0	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene		40.6	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene		46.2	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	48.5	50.0	ug/L	97	(71%-134%)
Bromofluorobenzene	49.4	50.0	ug/L	99	(70%-131%)
Toluene-d8	49.7	50.0	ug/L	99	(74%-124%)

Volatile
Certificate of Analysis
Sample Summary

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 1203985582

Client Sample: QC for batch 1744716

Client ID: LCS for batch 1744716

Batch ID: 1744716

Run Date: 03/08/2018 00:41

Prep Date: 03/08/2018 00:41

Data File: 030718V9\9I331L.D

Client: ARSL004

Method: SW-846:8260B

Inst: VOA9.I

Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC

SOP Ref: GL-OA-E-038

Dilution: 1

Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
630-20-6	1,1,1,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
71-55-6	1,1,1-Trichloroethane	U	0.300	ug/L	0.300	1.00
79-34-5	1,1,2,2-Tetrachloroethane	U	0.300	ug/L	0.300	1.00
79-00-5	1,1,2-Trichloroethane	U	0.300	ug/L	0.300	1.00
75-34-3	1,1-Dichloroethane	U	0.300	ug/L	0.300	1.00
75-35-4	1,1-Dichloroethylene	U	0.300	ug/L	0.300	1.00
563-58-6	1,1-Dichloropropene	U	0.300	ug/L	0.300	1.00
87-61-6	1,2,3-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
96-18-4	1,2,3-Trichloropropane	U	0.300	ug/L	0.300	1.00
120-82-1	1,2,4-Trichlorobenzene	U	0.300	ug/L	0.300	1.00
95-63-6	1,2,4-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
96-12-8	1,2-Dibromo-3-chloropropane	U	0.500	ug/L	0.500	1.00
106-93-4	1,2-Dibromoethane	U	0.300	ug/L	0.300	1.00
95-50-1	1,2-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
107-06-2	1,2-Dichloroethane	U	0.300	ug/L	0.300	1.00
78-87-5	1,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
108-67-8	1,3,5-Trimethylbenzene	U	0.300	ug/L	0.300	1.00
541-73-1	1,3-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
142-28-9	1,3-Dichloropropane	U	0.300	ug/L	0.300	1.00
106-46-7	1,4-Dichlorobenzene	U	0.300	ug/L	0.300	1.00
594-20-7	2,2-Dichloropropane	U	0.300	ug/L	0.300	1.00
78-93-3	2-Butanone	U	1.50	ug/L	1.50	5.00
126-99-8	2-Chloro-1,3-butadiene		44.7	ug/L	0.300	1.00
95-49-8	2-Chlorotoluene	U	0.300	ug/L	0.300	1.00
591-78-6	2-Hexanone	U	1.50	ug/L	1.50	5.00
106-43-4	4-Chlorotoluene	U	0.300	ug/L	0.300	1.00
99-87-6	4-Isopropyltoluene	U	0.300	ug/L	0.300	1.00
108-10-1	4-Methyl-2-pentanone	U	1.50	ug/L	1.50	5.00
67-64-1	Acetone	U	1.50	ug/L	1.50	10.0
75-05-8	Acetonitrile	U	8.00	ug/L	8.00	25.0
107-02-8	Acrolein		236	ug/L	1.50	5.00
107-13-1	Acrylonitrile		234	ug/L	1.50	5.00
107-05-1	Allyl chloride		214	ug/L	1.50	5.00
71-43-2	Benzene	U	0.300	ug/L	0.300	1.00
108-86-1	Bromobenzene	U	0.300	ug/L	0.300	1.00
74-97-5	Bromochloromethane	U	0.300	ug/L	0.300	1.00
75-27-4	Bromodichloromethane	U	0.300	ug/L	0.300	1.00
75-25-2	Bromoform	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

SDG Number: 2018-1811
Lab Sample ID: 1203985582
Client Sample: QC for batch 1744716
Client ID: LCS for batch 1744716
Batch ID: 1744716
Run Date: 03/08/2018 00:41
Prep Date: 03/08/2018 00:41
Data File: 030718V9\9I331L.D

Client: ARSL004
Method: SW-846:8260B
Inst: VOA9.I
Analyst: RXY1

Column: DB-624

Matrix: WATER

Project: QC
SOP Ref: GL-OA-E-038
Dilution: 1
Purge Vol: 5 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
74-83-9	Bromomethane	U	0.300	ug/L	0.300	1.00
75-15-0	Carbon disulfide	U	1.50	ug/L	1.50	5.00
56-23-5	Carbon tetrachloride	U	0.300	ug/L	0.300	1.00
108-90-7	Chlorobenzene	U	0.300	ug/L	0.300	1.00
75-00-3	Chloroethane	U	0.300	ug/L	0.300	1.00
67-66-3	Chloroform	U	0.300	ug/L	0.300	1.00
74-87-3	Chloromethane	U	0.300	ug/L	0.300	1.00
124-48-1	Dibromochloromethane	U	0.300	ug/L	0.300	1.00
74-95-3	Dibromomethane	U	0.300	ug/L	0.300	1.00
75-71-8	Dichlorodifluoromethane	U	0.300	ug/L	0.300	1.00
60-29-7	Ethyl ether	U	0.300	ug/L	0.300	1.00
97-63-2	Ethyl methacrylate		225	ug/L	1.50	5.00
100-41-4	Ethylbenzene	U	0.300	ug/L	0.300	1.00
87-68-3	Hexachlorobutadiene	U	0.300	ug/L	0.300	1.00
74-88-4	Iodomethane	U	1.50	ug/L	1.50	5.00
78-83-1	Isobutyl alcohol		2330	ug/L	15.0	50.0
98-82-8	Isopropylbenzene	U	0.300	ug/L	0.300	1.00
126-98-7	Methacrylonitrile		226	ug/L	1.50	5.00
80-62-6	Methyl methacrylate		232	ug/L	1.50	5.00
75-09-2	Methylene chloride	U	1.00	ug/L	1.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
107-12-0	Propionitrile		227	ug/L	1.50	5.00
100-42-5	Styrene	U	0.300	ug/L	0.300	1.00
127-18-4	Tetrachloroethylene	U	0.300	ug/L	0.300	1.00
108-88-3	Toluene	U	0.300	ug/L	0.300	1.00
79-01-6	Trichloroethylene	U	0.300	ug/L	0.300	1.00
75-69-4	Trichlorofluoromethane	U	0.300	ug/L	0.300	1.00
76-13-1	Trichlorotrifluoroethane		211	ug/L	2.00	5.00
108-05-4	Vinyl acetate	U	1.50	ug/L	1.50	5.00
75-01-4	Vinyl chloride	U	0.300	ug/L	0.300	1.00
156-59-2	cis-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-01-5	cis-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00
179601-23-1	m,p-Xylenes	U	0.300	ug/L	0.300	2.00
71-36-3	n-Butyl alcohol	U	15.0	ug/L	15.0	50.0
104-51-8	n-Butylbenzene	U	0.300	ug/L	0.300	1.00
103-65-1	n-Propylbenzene	U	0.300	ug/L	0.300	1.00
95-47-6	o-Xylene	U	0.300	ug/L	0.300	1.00
135-98-8	sec-Butylbenzene	U	0.300	ug/L	0.300	1.00

Volatile
Certificate of Analysis
Sample Summary

SDG Number:	2018-1811	Matrix:	WATER
Lab Sample ID:	1203985582		
Client Sample:	QC for batch 1744716	Client:	ARSL004
Client ID:	LCS for batch 1744716	Method:	SW-846:8260B
Batch ID:	1744716	Inst:	VOA9.I
Run Date:	03/08/2018 00:41	Analyst:	RXY1
Prep Date:	03/08/2018 00:41		
Data File:	030718V9\9I331L.D	Column:	DB-624

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
1634-04-4	tert-Butyl methyl ether	U	0.300	ug/L	0.300	1.00
98-06-6	tert-Butylbenzene	U	0.300	ug/L	0.300	1.00
156-60-5	trans-1,2-Dichloroethylene	U	0.300	ug/L	0.300	1.00
10061-02-6	trans-1,3-Dichloropropylene	U	0.300	ug/L	0.300	1.00

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
1,2-Dichloroethane-d4	50.2	50.0	ug/L 100	(71%-134%)
Bromofluorobenzene	48.4	50.0	ug/L 97	(70%-131%)
Toluene-d8	51.7	50.0	ug/L 103	(74%-124%)

Semi-Volatile Analysis

Case Narrative

**GC/MS Semivolatile
Technical Case Narrative
ARS International, LLC (ARSL)
SDG #: 2018-1811
Work Order #: 444566**

Method/Analysis Information

Procedure:	Analysis of Semivolatile Organic Compounds by Gas Chromatography/Mass Spectrometry
Analytical Method:	SW846 3510C/8270D
Prep Method:	SW846 3510C
Analytical Batch Number:	1742249
Prep Batch Number:	1742247

Sample Analysis

The following samples were analyzed using the analytical protocol as established in SW846 3510C/8270D:

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566009	CAWA-18-54
444566012	CAWA-18-56
1203978892	Method Blank (MB)
1203978893	Laboratory Control Sample (LCS)
1203978894	444566012(CAWA-18-56) Matrix Spike (MS)
1203978895	444566012(CAWA-18-56) Matrix Spike Duplicate (MSD)

The samples in this SDG were analyzed on an "as received" basis.

Preparation/Analytical Method Verification

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-OA-E-009 REV# 40.

Raw data reports are processed and reviewed by the analyst using the data analysis software package. False positives have been removed from the quantitation reports per standard operating procedures (SOP).

Calibration Information

A complete list of the initial calibration data files are shown in the Calibration History report located in the Standard Data section of the data package. The various calibration mixes may not be calibrated using all of the calibration levels. In addition, not all of the mixes are calibrated using the same levels.

Diphenylamine has now superseded N-Nitroso-diphenylamine on Quantitation Reports, Initial Calibration Reports, Calibration Check Standard Reports, etc. Previous versions of EPA Methodologies referenced N-Nitroso-diphenylamine. However, as stated in EPA Methodology, "N-Nitroso-diphenylamine decomposes in the gas chromatographic inlet and cannot be separated from Diphenylamine." Studies of these two compounds at GEL, both independent of each other and together, showed that they not only co-elute, but also have similar mass spectra. N-Nitroso-diphenylamine and Diphenylamine will be reported as Diphenylamine on all reports and forms.

Initial Calibration

All initial calibration requirements have been met for this sample delivery group (SDG) in this batch. A second source initial calibration verification (ICV) was included in the standard section directly behind the initial calibration.

CCV Requirements

All Calibration Verification Standards (CCV) did not meet the acceptance criteria as outlined in Method 8270D for samples 444566002 (CAPA-18-2), 444566006 (CAPA-18-12) and 444566012 (CAWA-18-56) and the associated QC. However, the method allows for a designated number of outliers dependent on the requested analyte list. This SDG satisfied the 8270D outlier acceptance criteria. If required, a CRDL was analyzed after the CCVs to demonstrate that there is adequate sensitivity to detect the failed compounds at the applicable lower quantitation limit.

Quality Control (QC) Information

Method Blank (MB) Statement

The MB analyzed with this SDG in this batch met the acceptance criteria.

Surrogate Recoveries

Sample (See Below) did not meet surrogate recovery acceptance criteria. Since there is insufficient sample remaining to perform a re-extraction, the results from this extraction are reported.

Sample	Analyte	Value
444566009 (CAWA-18-54)	2, 4, 6-Tribromophenol	24* (32%-124%)
	2-Fluorobiphenyl	21* (32%-112%)
	Nitrobenzene-d5	24* (36%-115%)
	Phenol-d5	12* (15%-91%)
	p-Terphenyl-d14	25* (36%-121%)

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries met the acceptance limits.

QC Sample Designation

Sample 444566012 (CAWA-18-56) was selected for analysis as the matrix spike and matrix spike duplicate.

Spike Recovery Statement

The MS and MSD recoveries were within the established acceptance limits.

MS/MSD Relative Percent Difference (RPD) Statement

The RPD values between the MS and MSD met the acceptance limits.

Internal Standard (ISTD) Acceptance

Sample 444566002 (CAPA-18-2) failed ISTD acceptance criteria. The sample was re-analyzed and confirmed the failures. The initial analysis data are reported. The re-analysis raw data have been placed in the Miscellaneous Section where applicable.

Technical Information:**Holding Time Specifications**

All samples in this SDG in this batch met the specified holding time. GEL assigns holding times based on the associated methodology that assigns the date and time from sample collection or sample receipt. Those holding times expressed in hours are calculated in the ALPHALIMS system. Those holding times expressed as days expire at midnight on the day of expiration.

Preparation/Analytical Method Verification

All procedures were performed as stated in the SOP. All reported compound mass spectra met the detection specifications in the method.

Sample Dilutions

The samples in this SDG in this batch did not require dilutions.

Sample Re-extraction/Re-analysis

Re-extractions or re-analyses were not required in this SDG in this analytical batch unless confirmations or dilutions were required.

Miscellaneous Information:**Manual Integrations**

Manual integrations were not required on samples in this batch in this SDG for detected target analytes or surrogates.

TIC Comment

Tentatively identified compounds (TIC) were requested for samples 444566002 (CAPA-18-2), 444566006 (CAPA-18-12), 444566009 (CAWA-18-54) and 444566012 (CAWA-18-56) in this SDG in this batch.

Additional Comments

Additional comments were not required for the SDG associated samples in this batch.

Electronic Package Comment

The following package was generated using an electronic data processing program referred to as "virtual packaging". In an effort to increase quality and efficiency, the laboratory is developing systems to eventually generate all data packages electronically. The following change from "traditional" packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. An electronic signature page inserted after the case narrative of each electronic package will indicate the reviewer name associated with the generation of the data and package. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

System Configuration

The Semi-Volatile-GC/MS analysis was performed on the following instrument configuration:

Instrument ID	Instrument	System Configuration	Column ID	Column Description
MSD3.I	Agilent 7890A/5975C GC/MS w/ 7683 Autosampler	HP7890A/HP5975C	DB-5MS	25m x 0.2mm, 0.33um (5% Phenylmethylpolysiloxane)

Certification Statement

Where the analytical method has been performed under NELAP certification, the analysis has met all of the requirements of the NELAC standard unless otherwise noted in the analytical case narrative.

GEL LABORATORIES LLC

2040 Savage Road Charleston SC 29407 - (843) 556-8171 - www.gel.com

Qualifier Definition Report for

ARSL004 ARS International, LLC (ARS-LANS-MTOA6-25093-GEL)

Client SDG: 2018-1811 GEL Work Order: 444566

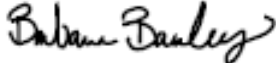
The Qualifiers in this report are defined as follows:

- * A quality control analyte recovery is outside of specified acceptance criteria
- ** Analyte is a surrogate compound
- U Analyte was analyzed for, but not detected above the MDL, MDA, MDC or LOD.
- DL Indicates that sample is diluted.
- RA Indicates that sample is re-analyzed without re-extraction.
- RE Indicates that sample is re-extracted.

Review/Validation

GEL requires all analytical data to be verified by a qualified data reviewer. In addition, all CLP-like deliverables receive a third level review of the fractional data package.

The following data validator verified the information presented in this data report:

Signature: 

Name: Barbara Bailey

Date: 20 MAR 2018

Title: Data Validator

Sample Data Summary

**Semi-Volatile
Certificate of Analysis
Sample Summary**

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566002

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW846 3510C/8270D

SOP Ref: GL-OA-E-009

Inst: MSD3.I

Dilution: 1

Batch ID: 1742249

Analyst: JLD1

Inj. Vol: 1 uL

Run Date: 02/27/2018 22:32

Aliquot: 940 mL

Final Volume: 1 mL

Prep Date: 02/27/2018 08:00

Column: DB-5ms

Data File: s022718.s\3b2722.D

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
95-94-3	1,2,4,5-Tetrachlorobenzene	U	3.19	ug/L	3.19	10.6
120-82-1	1,2,4-Trichlorobenzene	U	3.19	ug/L	3.19	10.6
95-50-1	1,2-Dichlorobenzene	U	3.19	ug/L	3.19	10.6
122-66-7	Azobenzene	U	3.19	ug/L	3.19	10.6
	<i>1,2-Diphenylhydrazine</i>					
541-73-1	1,3-Dichlorobenzene	U	3.19	ug/L	3.19	10.6
106-46-7	1,4-Dichlorobenzene	U	3.19	ug/L	3.19	10.6
123-91-1	1,4-Dioxane	U	3.19	ug/L	3.19	10.6
90-12-0	1-Methylnaphthalene	U	0.319	ug/L	0.319	1.06
58-90-2	2,3,4,6-Tetrachlorophenol	U	3.19	ug/L	3.19	10.6
95-95-4	2,4,5-Trichlorophenol	U	3.19	ug/L	3.19	10.6
88-06-2	2,4,6-Trichlorophenol	U	3.19	ug/L	3.19	10.6
120-83-2	2,4-Dichlorophenol	U	3.19	ug/L	3.19	10.6
105-67-9	2,4-Dimethylphenol	U	3.19	ug/L	3.19	10.6
51-28-5	2,4-Dinitrophenol	U	5.32	ug/L	5.32	21.3
121-14-2	2,4-Dinitrotoluene	U	3.19	ug/L	3.19	10.6
606-20-2	2,6-Dinitrotoluene	U	3.19	ug/L	3.19	10.6
91-58-7	2-Chloronaphthalene	U	0.436	ug/L	0.436	1.06
95-57-8	2-Chlorophenol	U	3.19	ug/L	3.19	10.6
534-52-1	2-Methyl-4,6-dinitrophenol	U	3.19	ug/L	3.19	10.6
91-57-6	2-Methylnaphthalene	U	0.319	ug/L	0.319	1.06
88-75-5	2-Nitrophenol	U	3.19	ug/L	3.19	10.6
91-94-1	3,3'-Dichlorobenzidine	U	3.19	ug/L	3.19	10.6
101-55-3	4-Bromophenylphenylether	U	3.19	ug/L	3.19	10.6
59-50-7	Parachlorometa cresol	U	3.19	ug/L	3.19	10.6
	<i>4-Chloro-3-methylphenol</i>					
106-47-8	4-Chloroaniline	U	3.51	ug/L	3.51	10.6
7005-72-3	4-Chlorophenylphenylether	U	3.19	ug/L	3.19	10.6
100-02-7	4-Nitrophenol	U	3.19	ug/L	3.19	10.6
83-32-9	Acenaphthene	U	0.319	ug/L	0.319	1.06
208-96-8	Acenaphthylene	U	0.319	ug/L	0.319	1.06
62-53-3	Aniline	U	4.47	ug/L	4.47	10.6
120-12-7	Anthracene	U	0.319	ug/L	0.319	1.06
1912-24-9	Atrazine	U	3.19	ug/L	3.19	10.6
92-87-5	Benzidine	U	4.15	ug/L	4.15	10.6
56-55-3	Benzo(a)anthracene	U	0.319	ug/L	0.319	1.06
50-32-8	Benzo(a)pyrene	U	0.319	ug/L	0.319	1.06
205-99-2	Benzo(b)fluoranthene	U	0.319	ug/L	0.319	1.06
191-24-2	Benzo(ghi)perylene	U	0.319	ug/L	0.319	1.06

**Semi-Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811

Lab Sample ID: 444566002

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW846 3510C/8270D

SOP Ref: GL-OA-E-009

Inst: MSD3.I

Dilution: 1

Batch ID: 1742249

Analyst: JLD1

Inj. Vol: 1 uL

Run Date: 02/27/2018 22:32

Prep Date: 02/27/2018 08:00

Aliquot: 940 mL

Final Volume: 1 mL

Data File: s022718.s\s3b2722.D

Column: DB-5ms

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
207-08-9	Benzo(k)fluoranthene	U	0.319	ug/L	0.319	1.06
65-85-0	Benzoic acid	U	6.38	ug/L	6.38	21.3
100-51-6	Benzyl alcohol	U	3.19	ug/L	3.19	10.6
85-68-7	Butylbenzylphthalate	U	3.19	ug/L	3.19	10.6
218-01-9	Chrysene	U	0.319	ug/L	0.319	1.06
84-74-2	Di-n-butylphthalate	U	3.19	ug/L	3.19	10.6
117-84-0	Di-n-octylphthalate	U	3.19	ug/L	3.19	10.6
53-70-3	Dibenzo(a,h)anthracene	U	0.319	ug/L	0.319	1.06
132-64-9	Dibenzofuran	U	3.19	ug/L	3.19	10.6
84-66-2	Diethylphthalate	U	3.19	ug/L	3.19	10.6
131-11-3	Dimethylphthalate	U	3.19	ug/L	3.19	10.6
88-85-7	Dinoseb	U	3.19	ug/L	3.19	10.6
122-39-4	Diphenylamine	U	3.19	ug/L	3.19	10.6
206-44-0	Fluoranthene	U	0.319	ug/L	0.319	1.06
86-73-7	Fluorene	U	0.319	ug/L	0.319	1.06
118-74-1	Hexachlorobenzene	U	3.19	ug/L	3.19	10.6
87-68-3	Hexachlorobutadiene	U	3.19	ug/L	3.19	10.6
77-47-4	Hexachlorocyclopentadiene	U	3.19	ug/L	3.19	10.6
67-72-1	Hexachloroethane	U	3.19	ug/L	3.19	10.6
193-39-5	Indeno(1,2,3-cd)pyrene	U	0.319	ug/L	0.319	1.06
78-59-1	Isophorone	U	3.72	ug/L	3.72	10.6
62-75-9	N-Methyl-N-nitrosomethylamine	U	3.19	ug/L	3.19	10.6
924-16-3	N-Nitrosodi-n-butylamine	U	3.19	ug/L	3.19	10.6
55-18-5	N-Nitrosodiethylamine	U	3.19	ug/L	3.19	10.6
621-64-7	N-Nitrosodi--n-propylamine	U	3.19	ug/L	3.19	10.6
	<i>N-Nitrosodipropylamine</i>					
930-55-2	N-Nitrosopyrrolidine	U	3.19	ug/L	3.19	10.6
91-20-3	Naphthalene	U	0.319	ug/L	0.319	1.06
98-95-3	Nitrobenzene	U	3.19	ug/L	3.19	10.6
608-93-5	Pentachlorobenzene	U	3.19	ug/L	3.19	10.6
87-86-5	Pentachlorophenol	U	3.19	ug/L	3.19	10.6
85-01-8	Phenanthrene	U	0.319	ug/L	0.319	1.06
108-95-2	Phenol	U	3.19	ug/L	3.19	10.6
129-00-0	Pyrene	U	0.319	ug/L	0.319	1.06
110-86-1	Pyridine	U	3.19	ug/L	3.19	10.6
108-60-1	bis(2-Chloro-1-methylethyl)ether	U	3.19	ug/L	3.19	10.6
111-91-1	bis(2-Chloroethoxy)methane	U	3.19	ug/L	3.19	10.6
111-44-4	bis(2-Chloroethyl) ether	U	3.19	ug/L	3.19	10.6
117-81-7	bis(2-Ethylhexyl)phthalate	U	3.19	ug/L	3.19	1.06

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811
Lab Sample ID: 444566002

Client ID: CAPA-18-2
Batch ID: 1742249
Run Date: 02/27/2018 22:32
Prep Date: 02/27/2018 08:00
Data File: s022718.s\s3b2722.D

Date Collected: 02/21/2018 11:47
Date Received: 02/23/2018 09:25
Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 940 mL
Column: DB-5ms

Matrix: W

Project: ESHL00114
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
65794-96-9	m,p-Cresols	U	3.94	ug/L	3.94	10.6
99-09-2	3-Nitroaniline	U	3.19	ug/L	3.19	10.6
	<i>m-Nitroaniline</i>					
95-48-7	o-Cresol	U	3.19	ug/L	3.19	10.6
88-74-4	2-Nitroaniline	U	3.19	ug/L	3.19	10.6
	<i>o-Nitroaniline</i>					
100-01-6	4-Nitroaniline	U	3.19	ug/L	3.19	10.6
	<i>p-Nitroaniline</i>					

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
2,4,6-Tribromophenol	71.8	106	ug/L	68	(32%-124%)
2-Fluorobiphenyl	29.8	53.2	ug/L	56	(32%-112%)
2-Fluorophenol	44.4	106	ug/L	42	(15%-88%)
Nitrobenzene-d5	34.7	53.2	ug/L	65	(36%-115%)
Phenol-d5	28.5	106	ug/L	27	(15%-91%)
p-Terphenyl-d14	36.5	53.2	ug/L	69	(36%-121%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
No Tentatively Identified Compounds Found				ug/L		

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811

Lab Sample ID: 444566006

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW846 3510C/8270D

SOP Ref: GL-OA-E-009

Inst: MSD3.I

Dilution: 1

Batch ID: 1742249

Analyst: JLD1

Inj. Vol: 1 uL

Run Date: 02/27/2018 23:02

Aliquot: 890 mL

Final Volume: 1 mL

Prep Date: 02/27/2018 08:00

Column: DB-5ms

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
95-94-3	1,2,4,5-Tetrachlorobenzene	U	3.37	ug/L	3.37	11.2
120-82-1	1,2,4-Trichlorobenzene	U	3.37	ug/L	3.37	11.2
95-50-1	1,2-Dichlorobenzene	U	3.37	ug/L	3.37	11.2
122-66-7	Azobenzene	U	3.37	ug/L	3.37	11.2
	<i>1,2-Diphenylhydrazine</i>					
541-73-1	1,3-Dichlorobenzene	U	3.37	ug/L	3.37	11.2
106-46-7	1,4-Dichlorobenzene	U	3.37	ug/L	3.37	11.2
123-91-1	1,4-Dioxane	U	3.37	ug/L	3.37	11.2
90-12-0	1-Methylnaphthalene	U	0.337	ug/L	0.337	1.12
58-90-2	2,3,4,6-Tetrachlorophenol	U	3.37	ug/L	3.37	11.2
95-95-4	2,4,5-Trichlorophenol	U	3.37	ug/L	3.37	11.2
88-06-2	2,4,6-Trichlorophenol	U	3.37	ug/L	3.37	11.2
120-83-2	2,4-Dichlorophenol	U	3.37	ug/L	3.37	11.2
105-67-9	2,4-Dimethylphenol	U	3.37	ug/L	3.37	11.2
51-28-5	2,4-Dinitrophenol	U	5.62	ug/L	5.62	22.5
121-14-2	2,4-Dinitrotoluene	U	3.37	ug/L	3.37	11.2
606-20-2	2,6-Dinitrotoluene	U	3.37	ug/L	3.37	11.2
91-58-7	2-Chloronaphthalene	U	0.461	ug/L	0.461	1.12
95-57-8	2-Chlorophenol	U	3.37	ug/L	3.37	11.2
534-52-1	2-Methyl-4,6-dinitrophenol	U	3.37	ug/L	3.37	11.2
91-57-6	2-Methylnaphthalene	U	0.337	ug/L	0.337	1.12
88-75-5	2-Nitrophenol	U	3.37	ug/L	3.37	11.2
91-94-1	3,3'-Dichlorobenzidine	U	3.37	ug/L	3.37	11.2
101-55-3	4-Bromophenylphenylether	U	3.37	ug/L	3.37	11.2
59-50-7	Parachlorometa cresol	U	3.37	ug/L	3.37	11.2
	<i>4-Chloro-3-methylphenol</i>					
106-47-8	4-Chloroaniline	U	3.71	ug/L	3.71	11.2
7005-72-3	4-Chlorophenylphenylether	U	3.37	ug/L	3.37	11.2
100-02-7	4-Nitrophenol	U	3.37	ug/L	3.37	11.2
83-32-9	Acenaphthene	U	0.337	ug/L	0.337	1.12
208-96-8	Acenaphthylene	U	0.337	ug/L	0.337	1.12
62-53-3	Aniline	U	4.72	ug/L	4.72	11.2
120-12-7	Anthracene	U	0.337	ug/L	0.337	1.12
1912-24-9	Atrazine	U	3.37	ug/L	3.37	11.2
92-87-5	Benzidine	U	4.38	ug/L	4.38	11.2
56-55-3	Benzo(a)anthracene	U	0.337	ug/L	0.337	1.12
50-32-8	Benzo(a)pyrene	U	0.337	ug/L	0.337	1.12
205-99-2	Benzo(b)fluoranthene	U	0.337	ug/L	0.337	1.12
191-24-2	Benzo(ghi)perylene	U	0.337	ug/L	0.337	1.12

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811

Lab Sample ID: 444566006

Date Collected: 02/21/2018 11:47

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW846 3510C/8270D

SOP Ref: GL-OA-E-009

Inst: MSD3.I

Dilution: 1

Batch ID: 1742249

Analyst: JLD1

Inj. Vol: 1 uL

Run Date: 02/27/2018 23:02

Aliquot: 890 mL

Final Volume: 1 mL

Prep Date: 02/27/2018 08:00

Column: DB-5ms

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
207-08-9	Benzo(k)fluoranthene	U	0.337	ug/L	0.337	1.12
65-85-0	Benzoic acid	U	6.74	ug/L	6.74	22.5
100-51-6	Benzyl alcohol	U	3.37	ug/L	3.37	11.2
85-68-7	Butylbenzylphthalate	U	3.37	ug/L	3.37	11.2
218-01-9	Chrysene	U	0.337	ug/L	0.337	1.12
84-74-2	Di-n-butylphthalate	U	3.37	ug/L	3.37	11.2
117-84-0	Di-n-octylphthalate	U	3.37	ug/L	3.37	11.2
53-70-3	Dibenzo(a,h)anthracene	U	0.337	ug/L	0.337	1.12
132-64-9	Dibenzofuran	U	3.37	ug/L	3.37	11.2
84-66-2	Diethylphthalate	U	3.37	ug/L	3.37	11.2
131-11-3	Dimethylphthalate	U	3.37	ug/L	3.37	11.2
88-85-7	Dinoseb	U	3.37	ug/L	3.37	11.2
122-39-4	Diphenylamine	U	3.37	ug/L	3.37	11.2
206-44-0	Fluoranthene	U	0.337	ug/L	0.337	1.12
86-73-7	Fluorene	U	0.337	ug/L	0.337	1.12
118-74-1	Hexachlorobenzene	U	3.37	ug/L	3.37	11.2
87-68-3	Hexachlorobutadiene	U	3.37	ug/L	3.37	11.2
77-47-4	Hexachlorocyclopentadiene	U	3.37	ug/L	3.37	11.2
67-72-1	Hexachloroethane	U	3.37	ug/L	3.37	11.2
193-39-5	Indeno(1,2,3-cd)pyrene	U	0.337	ug/L	0.337	1.12
78-59-1	Isophorone	U	3.93	ug/L	3.93	11.2
62-75-9	N-Methyl-N-nitrosomethylamine	U	3.37	ug/L	3.37	11.2
924-16-3	N-Nitrosodi-n-butylamine	U	3.37	ug/L	3.37	11.2
55-18-5	N-Nitrosodiethylamine	U	3.37	ug/L	3.37	11.2
621-64-7	N-Nitrosodi-n-propylamine	U	3.37	ug/L	3.37	11.2
	<i>N-Nitrosodipropylamine</i>					
930-55-2	N-Nitrosopyrrolidine	U	3.37	ug/L	3.37	11.2
91-20-3	Naphthalene	U	0.337	ug/L	0.337	1.12
98-95-3	Nitrobenzene	U	3.37	ug/L	3.37	11.2
608-93-5	Pentachlorobenzene	U	3.37	ug/L	3.37	11.2
87-86-5	Pentachlorophenol	U	3.37	ug/L	3.37	11.2
85-01-8	Phenanthrene	U	0.337	ug/L	0.337	1.12
108-95-2	Phenol	U	3.37	ug/L	3.37	11.2
129-00-0	Pyrene	U	0.337	ug/L	0.337	1.12
110-86-1	Pyridine	U	3.37	ug/L	3.37	11.2
108-60-1	bis(2-Chloro-1-methylethyl)ether	U	3.37	ug/L	3.37	11.2
111-91-1	bis(2-Chloroethoxy)methane	U	3.37	ug/L	3.37	11.2
111-44-4	bis(2-Chloroethyl) ether	U	3.37	ug/L	3.37	11.2
117-81-7	bis(2-Ethylhexyl)phthalate	U	3.37	ug/L	3.37	1.12

**Semi-Volatile
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Sample Summary**

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SDG Number: 2018-1811
Lab Sample ID: 444566006

Client ID: CAPA-18-12
Batch ID: 1742249
Run Date: 02/27/2018 23:02
Prep Date: 02/27/2018 08:00
Data File: s022718.s\s3b2723.D

Date Collected: 02/21/2018 11:47
Date Received: 02/23/2018 09:25
Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 890 mL
Column: DB-5ms

Matrix: W

Project: ESHL00114
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
65794-96-9	m,p-Cresols	U	4.16	ug/L	4.16	11.2
99-09-2	3-Nitroaniline	U	3.37	ug/L	3.37	11.2
	<i>m</i> -Nitroaniline					
95-48-7	o-Cresol	U	3.37	ug/L	3.37	11.2
88-74-4	2-Nitroaniline	U	3.37	ug/L	3.37	11.2
	<i>o</i> -Nitroaniline					
100-01-6	4-Nitroaniline	U	3.37	ug/L	3.37	11.2
	<i>p</i> -Nitroaniline					

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
2,4,6-Tribromophenol	59.9	112	ug/L	53	(32%-124%)
2-Fluorobiphenyl	24.9	56.2	ug/L	44	(32%-112%)
2-Fluorophenol	34.6	112	ug/L	31	(15%-88%)
Nitrobenzene-d5	27.4	56.2	ug/L	49	(36%-115%)
Phenol-d5	22.7	112	ug/L	20	(15%-91%)
p-Terphenyl-d14	34.0	56.2	ug/L	61	(36%-121%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
No Tentatively Identified Compounds Found				ug/L		

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811

Lab Sample ID: 444566009

Date Collected: 02/21/2018 10:31

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW846 3510C/8270D

SOP Ref: GL-OA-E-009

Inst: MSD3.I

Dilution: 1

Batch ID: 1742249

Analyst: JLD1

Inj. Vol: 1 uL

Run Date: 02/27/2018 23:31

Prep Date: 02/27/2018 08:00

Aliquot: 970 mL

Final Volume: 1 mL

Data File: s022718.s\3b2724.D

Column: DB-5ms

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
95-94-3	1,2,4,5-Tetrachlorobenzene	U	3.09	ug/L	3.09	10.3
120-82-1	1,2,4-Trichlorobenzene	U	3.09	ug/L	3.09	10.3
95-50-1	1,2-Dichlorobenzene	U	3.09	ug/L	3.09	10.3
122-66-7	Azobenzene	U	3.09	ug/L	3.09	10.3
	<i>1,2-Diphenylhydrazine</i>					
541-73-1	1,3-Dichlorobenzene	U	3.09	ug/L	3.09	10.3
106-46-7	1,4-Dichlorobenzene	U	3.09	ug/L	3.09	10.3
123-91-1	1,4-Dioxane	U	3.09	ug/L	3.09	10.3
90-12-0	1-Methylnaphthalene	U	0.309	ug/L	0.309	1.03
58-90-2	2,3,4,6-Tetrachlorophenol	U	3.09	ug/L	3.09	10.3
95-95-4	2,4,5-Trichlorophenol	U	3.09	ug/L	3.09	10.3
88-06-2	2,4,6-Trichlorophenol	U	3.09	ug/L	3.09	10.3
120-83-2	2,4-Dichlorophenol	U	3.09	ug/L	3.09	10.3
105-67-9	2,4-Dimethylphenol	U	3.09	ug/L	3.09	10.3
51-28-5	2,4-Dinitrophenol	U	5.15	ug/L	5.15	20.6
121-14-2	2,4-Dinitrotoluene	U	3.09	ug/L	3.09	10.3
606-20-2	2,6-Dinitrotoluene	U	3.09	ug/L	3.09	10.3
91-58-7	2-Chloronaphthalene	U	0.423	ug/L	0.423	1.03
95-57-8	2-Chlorophenol	U	3.09	ug/L	3.09	10.3
534-52-1	2-Methyl-4,6-dinitrophenol	U	3.09	ug/L	3.09	10.3
91-57-6	2-Methylnaphthalene	U	0.309	ug/L	0.309	1.03
88-75-5	2-Nitrophenol	U	3.09	ug/L	3.09	10.3
91-94-1	3,3'-Dichlorobenzidine	U	3.09	ug/L	3.09	10.3
101-55-3	4-Bromophenylphenylether	U	3.09	ug/L	3.09	10.3
59-50-7	Parachlorometa cresol	U	3.09	ug/L	3.09	10.3
	<i>4-Chloro-3-methylphenol</i>					
106-47-8	4-Chloroaniline	U	3.40	ug/L	3.40	10.3
7005-72-3	4-Chlorophenylphenylether	U	3.09	ug/L	3.09	10.3
100-02-7	4-Nitrophenol	U	3.09	ug/L	3.09	10.3
83-32-9	Acenaphthene	U	0.309	ug/L	0.309	1.03
208-96-8	Acenaphthylene	U	0.309	ug/L	0.309	1.03
62-53-3	Aniline	U	4.33	ug/L	4.33	10.3
120-12-7	Anthracene	U	0.309	ug/L	0.309	1.03
1912-24-9	Atrazine	U	3.09	ug/L	3.09	10.3
92-87-5	Benzidine	U	4.02	ug/L	4.02	10.3
56-55-3	Benzo(a)anthracene	U	0.309	ug/L	0.309	1.03
50-32-8	Benzo(a)pyrene	U	0.309	ug/L	0.309	1.03
205-99-2	Benzo(b)fluoranthene	U	0.309	ug/L	0.309	1.03
191-24-2	Benzo(ghi)perylene	U	0.309	ug/L	0.309	1.03

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811

Lab Sample ID: 444566009

Date Collected: 02/21/2018 10:31

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW846 3510C/8270D

SOP Ref: GL-OA-E-009

Inst: MSD3.I

Dilution: 1

Batch ID: 1742249

Analyst: JLD1

Inj. Vol: 1 uL

Run Date: 02/27/2018 23:31

Prep Date: 02/27/2018 08:00

Aliquot: 970 mL

Final Volume: 1 mL

Data File: s022718.s\s3b2724.D

Column: DB-5ms

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
207-08-9	Benzo(k)fluoranthene	U	0.309	ug/L	0.309	1.03
65-85-0	Benzoic acid	U	6.19	ug/L	6.19	20.6
100-51-6	Benzyl alcohol	U	3.09	ug/L	3.09	10.3
85-68-7	Butylbenzylphthalate	U	3.09	ug/L	3.09	10.3
218-01-9	Chrysene	U	0.309	ug/L	0.309	1.03
84-74-2	Di-n-butylphthalate	U	3.09	ug/L	3.09	10.3
117-84-0	Di-n-octylphthalate	U	3.09	ug/L	3.09	10.3
53-70-3	Dibenzo(a,h)anthracene	U	0.309	ug/L	0.309	1.03
132-64-9	Dibenzofuran	U	3.09	ug/L	3.09	10.3
84-66-2	Diethylphthalate	U	3.09	ug/L	3.09	10.3
131-11-3	Dimethylphthalate	U	3.09	ug/L	3.09	10.3
88-85-7	Dinoseb	U	3.09	ug/L	3.09	10.3
122-39-4	Diphenylamine	U	3.09	ug/L	3.09	10.3
206-44-0	Fluoranthene	U	0.309	ug/L	0.309	1.03
86-73-7	Fluorene	U	0.309	ug/L	0.309	1.03
118-74-1	Hexachlorobenzene	U	3.09	ug/L	3.09	10.3
87-68-3	Hexachlorobutadiene	U	3.09	ug/L	3.09	10.3
77-47-4	Hexachlorocyclopentadiene	U	3.09	ug/L	3.09	10.3
67-72-1	Hexachloroethane	U	3.09	ug/L	3.09	10.3
193-39-5	Indeno(1,2,3-cd)pyrene	U	0.309	ug/L	0.309	1.03
78-59-1	Isophorone	U	3.61	ug/L	3.61	10.3
62-75-9	N-Methyl-N-nitrosomethylamine	U	3.09	ug/L	3.09	10.3
924-16-3	N-Nitrosodi-n-butylamine	U	3.09	ug/L	3.09	10.3
55-18-5	N-Nitrosodiethylamine	U	3.09	ug/L	3.09	10.3
621-64-7	N-Nitrosodi-n-propylamine	U	3.09	ug/L	3.09	10.3
	<i>N-Nitrosodipropylamine</i>					
930-55-2	N-Nitrosopyrrolidine	U	3.09	ug/L	3.09	10.3
91-20-3	Naphthalene	U	0.309	ug/L	0.309	1.03
98-95-3	Nitrobenzene	U	3.09	ug/L	3.09	10.3
608-93-5	Pentachlorobenzene	U	3.09	ug/L	3.09	10.3
87-86-5	Pentachlorophenol	U	3.09	ug/L	3.09	10.3
85-01-8	Phenanthrene	U	0.309	ug/L	0.309	1.03
108-95-2	Phenol	U	3.09	ug/L	3.09	10.3
129-00-0	Pyrene	U	0.309	ug/L	0.309	1.03
110-86-1	Pyridine	U	3.09	ug/L	3.09	10.3
108-60-1	bis(2-Chloro-1-methylethyl)ether	U	3.09	ug/L	3.09	10.3
111-91-1	bis(2-Chloroethoxy)methane	U	3.09	ug/L	3.09	10.3
111-44-4	bis(2-Chloroethyl) ether	U	3.09	ug/L	3.09	10.3
117-81-7	bis(2-Ethylhexyl)phthalate	U	3.09	ug/L	3.09	1.03

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811

Lab Sample ID: 444566009

Date Collected: 02/21/2018 10:31

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Client ID: CAWA-18-54

Method: SW846 3510C/8270D

SOP Ref: GL-OA-E-009

Batch ID: 1742249

Inst: MSD3.I

Dilution: 1

Run Date: 02/27/2018 23:31

Analyst: JLD1

Inj. Vol: 1 uL

Prep Date: 02/27/2018 08:00

Aliquot: 970 mL

Final Volume: 1 mL

Data File: s022718.s\s3b2724.D

Column: DB-5ms

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
65794-96-9	m,p-Cresols	U	3.81	ug/L	3.81	10.3
99-09-2	3-Nitroaniline	U	3.09	ug/L	3.09	10.3
	<i>m</i> -Nitroaniline					
95-48-7	o-Cresol	U	3.09	ug/L	3.09	10.3
88-74-4	2-Nitroaniline	U	3.09	ug/L	3.09	10.3
	<i>o</i> -Nitroaniline					
100-01-6	4-Nitroaniline	U	3.09	ug/L	3.09	10.3
	<i>p</i> -Nitroaniline					

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
2,4,6-Tribromophenol	24.3	103	ug/L 24	* (32%-124%)
2-Fluorobiphenyl	10.8	51.5	ug/L 21	* (32%-112%)
2-Fluorophenol	18.3	103	ug/L 18	(15%-88%)
Nitrobenzene-d5	12.4	51.5	ug/L 24	* (36%-115%)
Phenol-d5	12.6	103	ug/L 12	* (15%-91%)
p-Terphenyl-d14	12.7	51.5	ug/L 25	* (36%-121%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
000103-23-1	Hexanedioic acid, bis(2-ethylhexyl	14.283	4.7	ug/L	95	NJ
1000309-37-9	Oxalic acid, isobutyl tetradecyl e	16.941	5.47	ug/L	91	NJ
1000314-61-4	Carbonic acid, heptadecyl isobutyl	18.931	4.97	ug/L	91	NJ

**Semi-Volatile
Certificate of Analysis
Sample Summary**

Page 1 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566012

Date Collected: 02/21/2018 10:44

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Project: ESHL00114

Method: SW846 3510C/8270D

SOP Ref: GL-OA-E-009

Inst: MSD3.I

Dilution: 1

Batch ID: 1742249

Analyst: JLD1

Inj. Vol: 1 uL

Run Date: 02/28/2018 00:01

Prep Date: 02/27/2018 08:00

Aliquot: 1000 mL

Final Volume: 1 mL

Data File: s022718.s\3b2725.D

Column: DB-5ms

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
95-94-3	1,2,4,5-Tetrachlorobenzene	U	3.00	ug/L	3.00	10.0
120-82-1	1,2,4-Trichlorobenzene	U	3.00	ug/L	3.00	10.0
95-50-1	1,2-Dichlorobenzene	U	3.00	ug/L	3.00	10.0
122-66-7	Azobenzene	U	3.00	ug/L	3.00	10.0
	<i>1,2-Diphenylhydrazine</i>					
541-73-1	1,3-Dichlorobenzene	U	3.00	ug/L	3.00	10.0
106-46-7	1,4-Dichlorobenzene	U	3.00	ug/L	3.00	10.0
123-91-1	1,4-Dioxane	U	3.00	ug/L	3.00	10.0
90-12-0	1-Methylnaphthalene	U	0.300	ug/L	0.300	1.00
58-90-2	2,3,4,6-Tetrachlorophenol	U	3.00	ug/L	3.00	10.0
95-95-4	2,4,5-Trichlorophenol	U	3.00	ug/L	3.00	10.0
88-06-2	2,4,6-Trichlorophenol	U	3.00	ug/L	3.00	10.0
120-83-2	2,4-Dichlorophenol	U	3.00	ug/L	3.00	10.0
105-67-9	2,4-Dimethylphenol	U	3.00	ug/L	3.00	10.0
51-28-5	2,4-Dinitrophenol	U	5.00	ug/L	5.00	20.0
121-14-2	2,4-Dinitrotoluene	U	3.00	ug/L	3.00	10.0
606-20-2	2,6-Dinitrotoluene	U	3.00	ug/L	3.00	10.0
91-58-7	2-Chloronaphthalene	U	0.410	ug/L	0.410	1.00
95-57-8	2-Chlorophenol	U	3.00	ug/L	3.00	10.0
534-52-1	2-Methyl-4,6-dinitrophenol	U	3.00	ug/L	3.00	10.0
91-57-6	2-Methylnaphthalene	U	0.300	ug/L	0.300	1.00
88-75-5	2-Nitrophenol	U	3.00	ug/L	3.00	10.0
91-94-1	3,3'-Dichlorobenzidine	U	3.00	ug/L	3.00	10.0
101-55-3	4-Bromophenylphenylether	U	3.00	ug/L	3.00	10.0
59-50-7	Parachlorometa cresol	U	3.00	ug/L	3.00	10.0
	<i>4-Chloro-3-methylphenol</i>					
106-47-8	4-Chloroaniline	U	3.30	ug/L	3.30	10.0
7005-72-3	4-Chlorophenylphenylether	U	3.00	ug/L	3.00	10.0
100-02-7	4-Nitrophenol	U	3.00	ug/L	3.00	10.0
83-32-9	Acenaphthene	U	0.300	ug/L	0.300	1.00
208-96-8	Acenaphthylene	U	0.300	ug/L	0.300	1.00
62-53-3	Aniline	U	4.20	ug/L	4.20	10.0
120-12-7	Anthracene	U	0.300	ug/L	0.300	1.00
1912-24-9	Atrazine	U	3.00	ug/L	3.00	10.0
92-87-5	Benzidine	U	3.90	ug/L	3.90	10.0
56-55-3	Benzo(a)anthracene	U	0.300	ug/L	0.300	1.00
50-32-8	Benzo(a)pyrene	U	0.300	ug/L	0.300	1.00
205-99-2	Benzo(b)fluoranthene	U	0.300	ug/L	0.300	1.00
191-24-2	Benzo(ghi)perylene	U	0.300	ug/L	0.300	1.00

**Semi-Volatile
Certificate of Analysis
Sample Summary**

Page 2 of 3

SDG Number: 2018-1811

Lab Sample ID: 444566012

Date Collected: 02/21/2018 10:44

Date Received: 02/23/2018 09:25

Matrix: W

Client: ARSL004

Method: SW846 3510C/8270D

Project: ESHL00114

SOP Ref: GL-OA-E-009

Client ID: CAWA-18-56

Batch ID: 1742249

Inst: MSD3.I

Dilution: 1

Run Date: 02/28/2018 00:01

Analyst: JLD1

Inj. Vol: 1 uL

Prep Date: 02/27/2018 08:00

Aliquot: 1000 mL

Final Volume: 1 mL

Data File: s022718.s\s3b2725.D

Column: DB-5ms

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
207-08-9	Benzo(k)fluoranthene	U	0.300	ug/L	0.300	1.00
65-85-0	Benzoic acid	U	6.00	ug/L	6.00	20.0
100-51-6	Benzyl alcohol	U	3.00	ug/L	3.00	10.0
85-68-7	Butylbenzylphthalate	U	3.00	ug/L	3.00	10.0
218-01-9	Chrysene	U	0.300	ug/L	0.300	1.00
84-74-2	Di-n-butylphthalate	U	3.00	ug/L	3.00	10.0
117-84-0	Di-n-octylphthalate	U	3.00	ug/L	3.00	10.0
53-70-3	Dibenzo(a,h)anthracene	U	0.300	ug/L	0.300	1.00
132-64-9	Dibenzofuran	U	3.00	ug/L	3.00	10.0
84-66-2	Diethylphthalate	U	3.00	ug/L	3.00	10.0
131-11-3	Dimethylphthalate	U	3.00	ug/L	3.00	10.0
88-85-7	Dinoseb	U	3.00	ug/L	3.00	10.0
122-39-4	Diphenylamine	U	3.00	ug/L	3.00	10.0
206-44-0	Fluoranthene	U	0.300	ug/L	0.300	1.00
86-73-7	Fluorene	U	0.300	ug/L	0.300	1.00
118-74-1	Hexachlorobenzene	U	3.00	ug/L	3.00	10.0
87-68-3	Hexachlorobutadiene	U	3.00	ug/L	3.00	10.0
77-47-4	Hexachlorocyclopentadiene	U	3.00	ug/L	3.00	10.0
67-72-1	Hexachloroethane	U	3.00	ug/L	3.00	10.0
193-39-5	Indeno(1,2,3-cd)pyrene	U	0.300	ug/L	0.300	1.00
78-59-1	Isophorone	U	3.50	ug/L	3.50	10.0
62-75-9	N-Methyl-N-nitrosomethylamine	U	3.00	ug/L	3.00	10.0
924-16-3	N-Nitrosodi-n-butylamine	U	3.00	ug/L	3.00	10.0
55-18-5	N-Nitrosodiethylamine	U	3.00	ug/L	3.00	10.0
621-64-7	N-Nitrosodi-n-propylamine	U	3.00	ug/L	3.00	10.0
	<i>N-Nitrosodipropylamine</i>					
930-55-2	N-Nitrosopyrrolidine	U	3.00	ug/L	3.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
98-95-3	Nitrobenzene	U	3.00	ug/L	3.00	10.0
608-93-5	Pentachlorobenzene	U	3.00	ug/L	3.00	10.0
87-86-5	Pentachlorophenol	U	3.00	ug/L	3.00	10.0
85-01-8	Phenanthrene	U	0.300	ug/L	0.300	1.00
108-95-2	Phenol	U	3.00	ug/L	3.00	10.0
129-00-0	Pyrene	U	0.300	ug/L	0.300	1.00
110-86-1	Pyridine	U	3.00	ug/L	3.00	10.0
108-60-1	bis(2-Chloro-1-methylethyl)ether	U	3.00	ug/L	3.00	10.0
111-91-1	bis(2-Chloroethoxy)methane	U	3.00	ug/L	3.00	10.0
111-44-4	bis(2-Chloroethyl) ether	U	3.00	ug/L	3.00	10.0
117-81-7	bis(2-Ethylhexyl)phthalate	U	3.00	ug/L	3.00	1.00

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811
Lab Sample ID: 444566012

Client ID: CAWA-18-56
Batch ID: 1742249
Run Date: 02/28/2018 00:01
Prep Date: 02/27/2018 08:00
Data File: s022718.s\s3b2725.D

Date Collected: 02/21/2018 10:44
Date Received: 02/23/2018 09:25
Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 1000 mL
Column: DB-5ms

Matrix: W

Project: ESHL00114
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
65794-96-9	m,p-Cresols	U	3.70	ug/L	3.70	10.0
99-09-2	3-Nitroaniline	U	3.00	ug/L	3.00	10.0
95-48-7	<i>m</i> -Nitroaniline o-Cresol	U	3.00	ug/L	3.00	10.0
88-74-4	2-Nitroaniline	U	3.00	ug/L	3.00	10.0
100-01-6	<i>o</i> -Nitroaniline 4-Nitroaniline <i>p</i> -Nitroaniline	U	3.00	ug/L	3.00	10.0

Surrogate/Tracer recovery	Result	Nominal	Recovery%	Acceptable Limits
2,4,6-Tribromophenol	52.1	100	ug/L	52 (32%-124%)
2-Fluorobiphenyl	23.6	50.0	ug/L	47 (32%-112%)
2-Fluorophenol	35.1	100	ug/L	35 (15%-88%)
Nitrobenzene-d5	28.2	50.0	ug/L	56 (36%-115%)
Phenol-d5	22.1	100	ug/L	22 (15%-91%)
p-Terphenyl-d14	33.6	50.0	ug/L	67 (36%-121%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
	unknown	2.05	4.69	ug/L	0	J

Quality Control Summary

Semi-Volatile
Surrogate Recovery Report

Page 1 of 1

SDG Number: 2018-1811

Matrix Type: LIQUID

Sample ID	Client ID	2FP %REC	PHL %REC	NBZ %REC	FBP %REC	TBP %REC	TPH %REC
1203978892	MB for batch 1742247	35	21	59	59	41	61
1203978893	LCS for batch 1742247	55	44	75	68	75	65
444566002	CAPA-18-2	42	27	65	56	68	69
444566006	CAPA-18-12	31	20	49	44	53	61
444566009	CAWA-18-54	18	12 *	24 *	21 *	24 *	25 *
444566012	CAWA-18-56	35	22	56	47	52	67
1203978894	CAWA-18-56MS	57	45	64	65	60	57
1203978895	CAWA-18-56MSD	55	43	67	62	64	58

Surrogate**Acceptance Limits**

2FP	= 2-Fluorophenol	(15%-88%)
PHL	= Phenol-d5	(15%-91%)
NBZ	= Nitrobenzene-d5	(36%-115%)
FBP	= 2-Fluorobiphenyl	(32%-112%)
TBP	= 2,4,6-Tribromophenol	(32%-124%)
TPH	= p-Terphenyl-d14	(36%-121%)

* Recovery outside Acceptance Limits

Column to be used to flag recovery values

D Sample Diluted

Semi-Volatile
Quality Control Summary
Spike Recovery Report

Page 1 of 4

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1742247

Matrix: WATER

Lab Sample ID 1203978893

Instrument: MSD3.I

Analysis Date: 02/27/2018 18:36

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
62-75-9	LCS N-Methyl-N-nitrosomethylamine	50.0	0.0	27.5	55	30-88
110-86-1	LCS Pyridine	50.0	0.0	31.4	63	27-89
62-53-3	LCS Aniline	50.0	0.0	39.9	80	49-112
108-95-2	LCS Phenol	50.0	0.0	21.4	43	16-82
111-44-4	LCS bis(2-Chloroethyl) ether	50.0	0.0	39.6	79	51-111
95-57-8	LCS 2-Chlorophenol	50.0	0.0	37.3	75	49-105
541-73-1	LCS 1,3-Dichlorobenzene	50.0	0.0	33.9	68	37-95
106-46-7	LCS 1,4-Dichlorobenzene	50.0	0.0	34.1	68	38-96
95-50-1	LCS 1,2-Dichlorobenzene	50.0	0.0	34.8	70	39-97
108-60-1	LCS bis(2-Chloro-1-methylethyl)ether	50.0	0.0	41.7	83	44-123
100-51-6	LCS Benzyl alcohol	50.0	0.0	35.2	70	44-102
95-48-7	LCS o-Cresol	50.0	0.0	34.1	68	41-101
65794-96-9	LCS m,p-Cresols	50.0	0.0	37.4	75	43-102
621-64-7	LCS N-Nitrosodi--n-propylamine <i>N-Nitrosodipropylamine</i>	50.0	0.0	43.6	87	54-115
67-72-1	LCS Hexachloroethane	50.0	0.0	33.3	67	36-96
98-95-3	LCS Nitrobenzene	50.0	0.0	37.0	74	53-115
78-59-1	LCS Isophorone	50.0	0.0	35.4	71	56-117
88-75-5	LCS 2-Nitrophenol	50.0	0.0	37.0	74	51-113
105-67-9	LCS 2,4-Dimethylphenol	50.0	0.0	32.4	65	51-104
111-91-1	LCS bis(2-Chloroethoxy)methane	50.0	0.0	37.7	75	55-114
120-83-2	LCS 2,4-Dichlorophenol	50.0	0.0	37.1	74	53-109
65-85-0	LCS Benzoic acid	100	0.0	40.5	40	21-74

Semi-Volatile
Quality Control Summary
Spike Recovery Report

Page 2 of 4

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1742247

Matrix: WATER

Lab Sample ID 1203978893

Instrument: MSD3.I

Analysis Date: 02/27/2018 18:36

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
106-47-8	LCS 4-Chloroaniline	50.0	0.0	49.7	99	65-136
87-68-3	LCS Hexachlorobutadiene	50.0	0.0	34.3	69	35-98
59-50-7	LCS Parachlorometa cresol 4-Chloro-3-methylphenol	50.0	0.0	36.1	72	55-115
91-57-6	LCS 2-Methylnaphthalene	50.0	0.0	32.6	65	42-103
91-20-3	LCS Naphthalene	50.0	0.0	34.2	68	44-102
90-12-0	LCS 1-Methylnaphthalene	50.0	0.0	33.8	68	45-108
77-47-4	LCS Hexachlorocyclopentadiene	50.0	0.0	27.8	56	34-89
88-06-2	LCS 2,4,6-Trichlorophenol	50.0	0.0	35.0	70	55-120
95-95-4	LCS 2,4,5-Trichlorophenol	50.0	0.0	33.7	67	55-116
91-58-7	LCS 2-Chloronaphthalene	50.0	0.0	34.1	68	44-107
88-74-4	LCS 2-Nitroaniline o-Nitroaniline	50.0	0.0	37.0	74	53-121
99-09-2	LCS 3-Nitroaniline m-Nitroaniline	50.0	0.0	42.6	85	61-139
131-11-3	LCS Dimethylphthalate	50.0	0.0	38.4	77	60-122
606-20-2	LCS 2,6-Dinitrotoluene	50.0	0.0	34.6	69	59-122
121-14-2	LCS 2,4-Dinitrotoluene	50.0	0.0	33.2	66	57-124
208-96-8	LCS Acenaphthylene	50.0	0.0	35.4	71	50-113
83-32-9	LCS Acenaphthene	50.0	0.0	38.4	77	49-112
51-28-5	LCS 2,4-Dinitrophenol	50.0	0.0	34.3	69	34-122
132-64-9	LCS Dibenzofuran	50.0	0.0	36.7	73	50-111
58-90-2	LCS 2,3,4,6-Tetrachlorophenol	50.0	0.0	32.0	64	54-122
84-66-2	LCS Diethylphthalate	50.0	0.0	38.9	78	57-122
100-02-7	LCS 4-Nitrophenol	50.0	0.0	17.6	35	15-137

Semi-Volatile
Quality Control Summary
Spike Recovery Report

Page 3 of 4

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1742247

Matrix: WATER

Lab Sample ID 1203978893

Instrument: MSD3.I

Analysis Date: 02/27/2018 18:36

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
86-73-7	LCS Fluorene	50.0	0.0	36.7	73	52-114
7005-72-3	LCS 4-Chlorophenylphenylether	50.0	0.0	38.6	77	52-121
100-01-6	LCS 4-Nitroaniline	50.0	0.0	35.9	72	44-137
	<i>p</i> -Nitroaniline					
534-52-1	LCS 2-Methyl-4,6-dinitrophenol	50.0	0.0	34.6	69	45-124
122-39-4	LCS Diphenylamine	50.0	0.0	37.4	75	55-113
122-66-7	LCS Azobenzene	50.0	0.0	38.0	76	53-115
	<i>1,2-Diphenylhydrazine</i>					
101-55-3	LCS 4-Bromophenylphenylether	50.0	0.0	38.1	76	54-116
118-74-1	LCS Hexachlorobenzene	50.0	0.0	37.3	75	54-115
87-86-5	LCS Pentachlorophenol	50.0	0.0	39.4	79	41-116
85-01-8	LCS Phenanthrene	50.0	0.0	36.9	74	55-110
120-12-7	LCS Anthracene	50.0	0.0	37.1	74	56-112
84-74-2	LCS Di-n-butylphthalate	50.0	0.0	42.7	85	57-123
206-44-0	LCS Fluoranthene	50.0	0.0	41.3	83	54-118
129-00-0	LCS Pyrene	50.0	0.0	28.2	56	49-121
85-68-7	LCS Butylbenzylphthalate	50.0	0.0	35.4	71	52-125
117-81-7	LCS bis(2-Ethylhexyl)phthalate	50.0	0.0	40.0	80	52-125
56-55-3	LCS Benzo(a)anthracene	50.0	0.0	36.2	72	57-112
218-01-9	LCS Chrysene	50.0	0.0	36.7	73	58-117
117-84-0	LCS Di-n-octylphthalate	50.0	0.0	39.1	78	50-129
205-99-2	LCS Benzo(b)fluoranthene	50.0	0.0	34.3	69	41-118
207-08-9	LCS Benzo(k)fluoranthene	50.0	0.0	35.6	71	42-121
50-32-8	LCS Benzo(a)pyrene	50.0	0.0	37.0	74	40-118

Semi-Volatile
Quality Control Summary
Spike Recovery Report

Page 4 of 4

SDG Number: 2018-1811

Sample Type: Laboratory Control Sample

Client ID: LCS for batch 1742247

Matrix: WATER

Lab Sample ID 1203978893

Instrument: MSD3.I

Analysis Date: 02/27/2018 18:36

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
193-39-5	LCS Indeno(1,2,3-cd)pyrene	50.0	0.0	43.8	88	34-125
53-70-3	LCS Dibenzo(a,h)anthracene	50.0	0.0	45.3	91	38-129
191-24-2	LCS Benzo(ghi)perylene	50.0	0.0	44.4	89	33-131
123-91-1	LCS 1,4-Dioxane	50.0	0.0	27.6	55	38-78
930-55-2	LCS N-Nitrosopyrrolidine	50.0	0.0	40.9	82	54-113
95-94-3	LCS 1,2,4,5-Tetrachlorobenzene	50.0	0.0	37.0	74	44-102
1912-24-9	LCS Atrazine	50.0	0.0	32.4	65	60-131
92-87-5	LCS Benzidine	100	0.0	88.9	89	20-144
91-94-1	LCS 3,3'-Dichlorobenzidine	50.0	0.0	44.1	88	43-127
120-82-1	LCS 1,2,4-Trichlorobenzene	50.0	0.0	34.2	68	39-99

Semi-Volatile
Quality Control Summary
Spike Recovery Report

Page 1 of 8

SDG Number: 2018-1811

Sample Type: Matrix Spike

Client ID: CAWA-18-56MS

Matrix: W

Lab Sample ID 1203978894

Instrument: MSD3.I

Analysis Date: 02/28/2018 00:30

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
62-75-9	MS N-Methyl-N-nitrosomethylamine	115	0.00 U	65.7	57	25-106
110-86-1	MS Pyridine	115	0.00 U	71.9	63	24-93
62-53-3	MS Aniline	115	0.00 U	78.6	68	37-113
108-95-2	MS Phenol	115	0.00 U	50.3	44	23-82
111-44-4	MS bis(2-Chloroethyl) ether	115	0.00 U	74.6	65	39-114
95-57-8	MS 2-Chlorophenol	115	0.00 U	72.1	63	37-108
541-73-1	MS 1,3-Dichlorobenzene	115	0.00 U	63.7	55	27-97
106-46-7	MS 1,4-Dichlorobenzene	115	0.00 U	62.7	55	28-97
95-50-1	MS 1,2-Dichlorobenzene	115	0.00 U	64.4	56	28-99
108-60-1	MS bis(2-Chloro-1-methylethyl)ether	115	0.00 U	78.0	68	32-127
100-51-6	MS Benzyl alcohol	115	0.00 U	70.7	61	37-116
95-48-7	MS o-Cresol	115	0.00 U	68.6	60	34-109
65794-96-9	MS m,p-Cresols	115	0.00 U	76.6	67	36-120
621-64-7	MS N-Nitrosodi--n-propylamine <i>N-Nitrosodipropylamine</i>	115	0.00 U	77.6	68	42-118
67-72-1	MS Hexachloroethane	115	0.00 U	60.4	53	29-94
98-95-3	MS Nitrobenzene	115	0.00 U	71.8	62	38-123
78-59-1	MS Isophorone	115	0.00 U	65.1	57	43-120
88-75-5	MS 2-Nitrophenol	115	0.00 U	70.3	61	39-115
105-67-9	MS 2,4-Dimethylphenol	115	0.00 U	59.1	51	39-107
111-91-1	MS bis(2-Chloroethoxy)methane	115	0.00 U	69.3	60	42-118
120-83-2	MS 2,4-Dichlorophenol	115	0.00 U	69.5	60	40-111
65-85-0	MS Benzoic acid	230	0.00 U	106	46	17-95

Semi-Volatile
Quality Control Summary
Spike Recovery Report

Page 2 of 8

SDG Number: 2018-1811

Sample Type: Matrix Spike

Client ID: CAWA-18-56MS

Matrix: W

Lab Sample ID 1203978894

Instrument: MSD3.I

Analysis Date: 02/28/2018 00:30

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
106-47-8	MS 4-Chloroaniline	115	0.00 U	91.8	80	44-138
87-68-3	MS Hexachlorobutadiene	115	0.00 U	63.4	55	26-98
59-50-7	MS Parachlorometa cresol 4-Chloro-3-methylphenol	115	0.00 U	64.9	56	41-122
91-57-6	MS 2-Methylnaphthalene	115	0.00 U	61.2	53	29-109
91-20-3	MS Naphthalene	115	0.00 U	67.3	59	31-108
90-12-0	MS 1-Methylnaphthalene	115	0.00 U	63.8	56	33-112
77-47-4	MS Hexachlorocyclopentadiene	115	0.00 U	61.4	53	26-79
88-06-2	MS 2,4,6-Trichlorophenol	115	0.00 U	71.1	62	39-124
95-95-4	MS 2,4,5-Trichlorophenol	115	0.00 U	66.7	58	42-120
91-58-7	MS 2-Chloronaphthalene	115	0.00 U	71.1	62	29-113
88-74-4	MS 2-Nitroaniline o-Nitroaniline	115	0.00 U	72.4	63	41-121
99-09-2	MS 3-Nitroaniline m-Nitroaniline	115	0.00 U	81.1	71	42-144
131-11-3	MS Dimethylphthalate	115	0.00 U	70.5	61	45-128
606-20-2	MS 2,6-Dinitrotoluene	115	0.00 U	63.3	55	46-124
121-14-2	MS 2,4-Dinitrotoluene	115	0.00 U	61.2	53	45-125
208-96-8	MS Acenaphthylene	115	0.00 U	70.9	62	35-120
83-32-9	MS Acenaphthene	115	0.00 U	75.4	66	35-117
51-28-5	MS 2,4-Dinitrophenol	115	0.00 U	72.9	63	27-122
132-64-9	MS Dibenzofuran	115	0.00 U	70.9	62	38-113
58-90-2	MS 2,3,4,6-Tetrachlorophenol	115	0.00 U	60.7	53	40-128
84-66-2	MS Diethylphthalate	115	0.00 U	70.3	61	43-127
100-02-7	MS 4-Nitrophenol	115	0.00 U	46.5	40	17-85

Semi-Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Matrix Spike

Client ID: CAWA-18-56MS

Matrix: W

Lab Sample ID 1203978894

Instrument: MSD3.I

Analysis Date: 02/28/2018 00:30

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
86-73-7	MS Fluorene	115	0.00 U	67.8	59	39-117
7005-72-3	MS 4-Chlorophenylphenylether	115	0.00 U	70.8	62	39-121
100-01-6	MS 4-Nitroaniline	115	0.00 U	71.4	62	30-133
534-52-1	MS 2-Methyl-4,6-dinitrophenol	115	0.00 U	69.3	60	32-126
122-39-4	MS Diphenylamine	115	0.00 U	70.7	61	37-118
122-66-7	MS Azobenzene	115	0.00 U	73.0	64	38-120
101-55-3	MS 4-Bromophenylphenylether	115	0.00 U	70.2	61	39-121
118-74-1	MS Hexachlorobenzene	115	0.00 U	70.2	61	40-118
87-86-5	MS Pentachlorophenol	115	0.00 U	79.0	69	35-121
85-01-8	MS Phenanthrene	115	0.00 U	70.0	61	40-115
120-12-7	MS Anthracene	115	0.00 U	69.6	61	38-120
84-74-2	MS Di-n-butylphthalate	115	0.00 U	83.9	73	41-128
206-44-0	MS Fluoranthene	115	0.00 U	79.8	69	41-119
129-00-0	MS Pyrene	115	0.00 U	56.2	49	35-128
85-68-7	MS Butylbenzylphthalate	115	0.00 U	69.9	61	40-129
117-81-7	MS bis(2-Ethylhexyl)phthalate	115	0.00 U	78.3	68	38-131
56-55-3	MS Benzo(a)anthracene	115	0.00 U	69.0	60	39-120
218-01-9	MS Chrysene	115	0.00 U	70.6	61	41-124
117-84-0	MS Di-n-octylphthalate	115	0.00 U	79.9	70	37-134
205-99-2	MS Benzo(b)fluoranthene	115	0.00 U	64.8	56	31-122
207-08-9	MS Benzo(k)fluoranthene	115	0.00 U	67.3	59	33-123
50-32-8	MS Benzo(a)pyrene	115	0.00 U	69.5	60	32-118

Semi-Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Matrix Spike

Client ID: CAWA-18-56MS

Matrix: W

Lab Sample ID 1203978894

Instrument: MSD3.I

Analysis Date: 02/28/2018 00:30

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits
193-39-5	MS Indeno(1,2,3-cd)pyrene	115	0.00 U	77.1	67	27-121
53-70-3	MS Dibenzo(a,h)anthracene	115	0.00 U	78.9	69	30-125
191-24-2	MS Benzo(ghi)perylene	115	0.00 U	77.8	68	24-126
123-91-1	MS 1,4-Dioxane	115	0.00 U	64.7	56	24-110
930-55-2	MS N-Nitrosopyrrolidine	115	0.00 U	75.7	66	47-119
95-94-3	MS 1,2,4,5-Tetrachlorobenzene	115	0.00 U	79.2	69	32-101
1912-24-9	MS Atrazine	115	0.00 U	76.1	66	42-129
92-87-5	MS Benzidine	230	0.00 U	222	96	15-130
91-94-1	MS 3,3'-Dichlorobenzidine	115	0.00 U	87.2	76	34-124
120-82-1	MS 1,2,4-Trichlorobenzene	115	0.00 U	65.1	57	26-102

Semi-Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Matrix Spike Duplicate

Client ID: CAWA-18-56MSD

Matrix: W

Lab Sample ID 1203978895

Instrument: MSD3.I

Analysis Date: 02/28/2018 01:00

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L		Spike Conc. ug/L	Recovery %	Acceptance Limits	RPD %	Acceptance Limits
62-75-9	MSD N-Methyl-N-nitrosomethylamine	115	0.00	U	62.8	55	25-106	4	0-30
110-86-1	MSD Pyridine	115	0.00	U	64.9	56	24-93	10	0-30
62-53-3	MSD Aniline	115	0.00	U	73.9	64	37-113	6	0-30
108-95-2	MSD Phenol	115	0.00	U	48.6	42	23-82	3	0-30
111-44-4	MSD bis(2-Chloroethyl) ether	115	0.00	U	77.8	68	39-114	4	0-30
95-57-8	MSD 2-Chlorophenol	115	0.00	U	74.3	65	37-108	3	0-30
541-73-1	MSD 1,3-Dichlorobenzene	115	0.00	U	63.4	55	27-97	0	0-30
106-46-7	MSD 1,4-Dichlorobenzene	115	0.00	U	63.9	56	28-97	2	0-30
95-50-1	MSD 1,2-Dichlorobenzene	115	0.00	U	65.8	57	28-99	2	0-30
108-60-1	MSD bis(2-Chloro-1-methylethyl)ether	115	0.00	U	81.8	71	32-127	5	0-30
100-51-6	MSD Benzyl alcohol	115	0.00	U	75.2	65	37-116	6	0-30
95-48-7	MSD o-Cresol	115	0.00	U	72.8	63	34-109	6	0-30
65794-96-9	MSD m,p-Cresols	115	0.00	U	83.0	72	36-120	8	0-30
621-64-7	MSD N-Nitrosodi--n-propylamine N-Nitrosodipropylamine	115	0.00	U	85.4	74	42-118	10	0-30
67-72-1	MSD Hexachloroethane	115	0.00	U	60.8	53	29-94	1	0-30
98-95-3	MSD Nitrobenzene	115	0.00	U	75.7	66	38-123	5	0-30
78-59-1	MSD Isophorone	115	0.00	U	69.9	61	43-120	7	0-30
88-75-5	MSD 2-Nitrophenol	115	0.00	U	75.4	66	39-115	7	0-30
105-67-9	MSD 2,4-Dimethylphenol	115	0.00	U	60.4	53	39-107	2	0-30
111-91-1	MSD bis(2-Chloroethoxy)methane	115	0.00	U	74.3	65	42-118	7	0-30
120-83-2	MSD 2,4-Dichlorophenol	115	0.00	U	75.2	65	40-111	8	0-30
65-85-0	MSD Benzoic acid	230	0.00	U	116	50	17-95	9	0-30

Semi-Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Matrix Spike Duplicate

Client ID: CAWA-18-56MSD

Matrix: W

Lab Sample ID 1203978895

Instrument: MSD3.I

Analysis Date: 02/28/2018 01:00

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits	RPD %	Acceptance Limits
106-47-8	MSD 4-Chloroaniline	115	0.00	U 90.5	79	44-138	1	0-30
87-68-3	MSD Hexachlorobutadiene	115	0.00	U 63.3	55	26-98	0	0-30
59-50-7	MSD Parachlorometa cresol 4-Chloro-3-methylphenol	115	0.00	U 76.0	66	41-122	16	0-30
91-57-6	MSD 2-Methylnaphthalene	115	0.00	U 68.3	59	29-109	11	0-30
91-20-3	MSD Naphthalene	115	0.00	U 70.7	61	31-108	5	0-30
90-12-0	MSD 1-Methylnaphthalene	115	0.00	U 70.4	61	33-112	10	0-30
77-47-4	MSD Hexachlorocyclopentadiene	115	0.00	U 54.7	48	26-79	12	0-30
88-06-2	MSD 2,4,6-Trichlorophenol	115	0.00	U 71.1	62	39-124	0	0-30
95-95-4	MSD 2,4,5-Trichlorophenol	115	0.00	U 69.3	60	42-120	4	0-30
91-58-7	MSD 2-Chloronaphthalene	115	0.00	U 69.7	61	29-113	2	0-30
88-74-4	MSD 2-Nitroaniline o-Nitroaniline	115	0.00	U 76.3	66	41-121	5	0-30
99-09-2	MSD 3-Nitroaniline m-Nitroaniline	115	0.00	U 82.4	72	42-144	2	0-30
131-11-3	MSD Dimethylphthalate	115	0.00	U 73.1	64	45-128	4	0-30
606-20-2	MSD 2,6-Dinitrotoluene	115	0.00	U 66.8	58	46-124	5	0-30
121-14-2	MSD 2,4-Dinitrotoluene	115	0.00	U 64.2	56	45-125	5	0-30
208-96-8	MSD Acenaphthylene	115	0.00	U 71.9	63	35-120	1	0-30
83-32-9	MSD Acenaphthene	115	0.00	U 75.5	66	35-117	0	0-30
51-28-5	MSD 2,4-Dinitrophenol	115	0.00	U 79.4	69	27-122	9	0-30
132-64-9	MSD Dibenzofuran	115	0.00	U 73.4	64	38-113	4	0-30
58-90-2	MSD 2,3,4,6-Tetrachlorophenol	115	0.00	U 64.0	56	40-128	5	0-30
84-66-2	MSD Diethylphthalate	115	0.00	U 72.6	63	43-127	3	0-30
100-02-7	MSD 4-Nitrophenol	115	0.00	U 45.5	40	17-85	2	0-30

Semi-Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Matrix Spike Duplicate

Client ID: CAWA-18-56MSD

Matrix: W

Lab Sample ID 1203978895

Instrument: MSD3.I

Analysis Date: 02/28/2018 01:00

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits	RPD %	Acceptance Limits
86-73-7	MSD Fluorene	115	0.00 U	71.8	62	39-117	6	0-30
7005-72-3	MSD 4-Chlorophenylphenylether	115	0.00 U	74.8	65	39-121	6	0-30
100-01-6	MSD 4-Nitroaniline <i>p</i> -Nitroaniline	115	0.00 U	75.3	66	30-133	5	0-30
534-52-1	MSD 2-Methyl-4,6-dinitrophenol	115	0.00 U	73.5	64	32-126	6	0-30
122-39-4	MSD Diphenylamine	115	0.00 U	74.3	65	37-118	5	0-30
122-66-7	MSD Azobenzene <i>1,2-Diphenylhydrazine</i>	115	0.00 U	76.5	67	38-120	5	0-30
101-55-3	MSD 4-Bromophenylphenylether	115	0.00 U	72.7	63	39-121	4	0-30
118-74-1	MSD Hexachlorobenzene	115	0.00 U	72.2	63	40-118	3	0-30
87-86-5	MSD Pentachlorophenol	115	0.00 U	85.0	74	35-121	7	0-30
85-01-8	MSD Phenanthrene	115	0.00 U	72.9	63	40-115	4	0-30
120-12-7	MSD Anthracene	115	0.00 U	72.1	63	38-120	4	0-30
84-74-2	MSD Di-n-butylphthalate	115	0.00 U	84.8	74	41-128	1	0-30
206-44-0	MSD Fluoranthene	115	0.00 U	81.0	70	41-119	1	0-30
129-00-0	MSD Pyrene	115	0.00 U	57.7	50	35-128	3	0-30
85-68-7	MSD Butylbenzylphthalate	115	0.00 U	71.8	62	40-129	3	0-30
117-81-7	MSD bis(2-Ethylhexyl)phthalate	115	0.00 U	81.1	71	38-131	3	0-30
56-55-3	MSD Benzo(a)anthracene	115	0.00 U	71.9	63	39-120	4	0-30
218-01-9	MSD Chrysene	115	0.00 U	73.6	64	41-124	4	0-30
117-84-0	MSD Di-n-octylphthalate	115	0.00 U	83.7	73	37-134	5	0-30
205-99-2	MSD Benzo(b)fluoranthene	115	0.00 U	68.9	60	31-122	6	0-30
207-08-9	MSD Benzo(k)fluoranthene	115	0.00 U	70.4	61	33-123	5	0-30
50-32-8	MSD Benzo(a)pyrene	115	0.00 U	73.7	64	32-118	6	0-30

Semi-Volatile
Quality Control Summary
Spike Recovery Report

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SDG Number: 2018-1811

Sample Type: Matrix Spike Duplicate

Client ID: CAWA-18-56MSD

Matrix: W

Lab Sample ID 1203978895

Instrument: MSD3.I

Analysis Date: 02/28/2018 01:00

Dilution: 1

Analyst: JLD1

Prep Batch ID: 1742247

Inj. Vol: 1 uL

Batch ID: 1742249

CAS No	Parmname	Amount Added ug/L	Sample Conc. ug/L	Spike Conc. ug/L	Recovery %	Acceptance Limits	RPD %	Acceptance Limits
193-39-5	MSD Indeno(1,2,3-cd)pyrene	115	0.00	U	76.3	66	27-121	1 0-30
53-70-3	MSD Dibenzo(a,h)anthracene	115	0.00	U	77.7	68	30-125	1 0-30
191-24-2	MSD Benzo(ghi)perylene	115	0.00	U	74.5	65	24-126	4 0-30
123-91-1	MSD 1,4-Dioxane	115	0.00	U	60.6	53	24-110	6 0-30
930-55-2	MSD N-Nitrosopyrrolidine	115	0.00	U	84.6	74	47-119	11 0-30
95-94-3	MSD 1,2,4,5-Tetrachlorobenzene	115	0.00	U	74.8	65	32-101	6 0-30
1912-24-9	MSD Atrazine	115	0.00	U	76.7	67	42-129	1 0-30
92-87-5	MSD Benzidine	230	0.00	U	202	88	15-130	9 0-30
91-94-1	MSD 3,3'-Dichlorobenzidine	115	0.00	U	91.1	79	34-124	4 0-30
120-82-1	MSD 1,2,4-Trichlorobenzene	115	0.00	U	67.3	59	26-102	3 0-30

Method Blank Summary

Page 1 of 1

SDG Number:	2018-1811	Client:	ARSL004	Matrix:	WATER
Client ID:	MB for batch 1742247	Instrument ID:	MSD3.I	Data File:	s022718.s\s3b2713.D
Lab Sample ID:	1203978892	Prep Date:	02/27/2018 08:00	Analyzed:	02/27/18 18:06
Column:	DB-5ms				

This method blank applies to the following samples and quality control samples:

Client Sample ID	Lab Sample ID	File ID	Date Analyzed	Time Analyzed
01 LCS for batch 1742247	1203978893	s022718.s\s3b2714.D	02/27/18	1836
02 CAPA-18-2	444566002	s022718.s\s3b2722.D	02/27/18	2232
03 CAPA-18-12	444566006	s022718.s\s3b2723.D	02/27/18	2302
04 CAWA-18-54	444566009	s022718.s\s3b2724.D	02/27/18	2331
05 CAWA-18-56	444566012	s022718.s\s3b2725.D	02/28/18	0001
06 CAWA-18-56MS	1203978894	s022718.s\s3b2726.D	02/28/18	0030
07 CAWA-18-56MSD	1203978895	s022718.s\s3b2727.D	02/28/18	0100

Quality Control Data

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811
Lab Sample ID: 1203978892
Client Sample: QC for batch 1742247
Client ID: MB for batch 1742247
Batch ID: 1742249
Run Date: 02/27/2018 18:06
Prep Date: 02/27/2018 08:00
Data File: s022718.s\3b2713.D

Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 1000 mL
Column: DB-5ms

Matrix: WATER
Project: QC
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
95-94-3	1,2,4,5-Tetrachlorobenzene	U	3.00	ug/L	3.00	10.0
120-82-1	1,2,4-Trichlorobenzene	U	3.00	ug/L	3.00	10.0
95-50-1	1,2-Dichlorobenzene	U	3.00	ug/L	3.00	10.0
122-66-7	Azobenzene	U	3.00	ug/L	3.00	10.0
	<i>1,2-Diphenylhydrazine</i>					
541-73-1	1,3-Dichlorobenzene	U	3.00	ug/L	3.00	10.0
106-46-7	1,4-Dichlorobenzene	U	3.00	ug/L	3.00	10.0
123-91-1	1,4-Dioxane	U	3.00	ug/L	3.00	10.0
90-12-0	1-Methylnaphthalene	U	0.300	ug/L	0.300	1.00
58-90-2	2,3,4,6-Tetrachlorophenol	U	3.00	ug/L	3.00	10.0
95-95-4	2,4,5-Trichlorophenol	U	3.00	ug/L	3.00	10.0
88-06-2	2,4,6-Trichlorophenol	U	3.00	ug/L	3.00	10.0
120-83-2	2,4-Dichlorophenol	U	3.00	ug/L	3.00	10.0
105-67-9	2,4-Dimethylphenol	U	3.00	ug/L	3.00	10.0
51-28-5	2,4-Dinitrophenol	U	5.00	ug/L	5.00	20.0
121-14-2	2,4-Dinitrotoluene	U	3.00	ug/L	3.00	10.0
606-20-2	2,6-Dinitrotoluene	U	3.00	ug/L	3.00	10.0
91-58-7	2-Chloronaphthalene	U	0.410	ug/L	0.410	1.00
95-57-8	2-Chlorophenol	U	3.00	ug/L	3.00	10.0
534-52-1	2-Methyl-4,6-dinitrophenol	U	3.00	ug/L	3.00	10.0
91-57-6	2-Methylnaphthalene	U	0.300	ug/L	0.300	1.00
88-75-5	2-Nitrophenol	U	3.00	ug/L	3.00	10.0
91-94-1	3,3'-Dichlorobenzidine	U	3.00	ug/L	3.00	10.0
101-55-3	4-Bromophenylphenylether	U	3.00	ug/L	3.00	10.0
59-50-7	Parachlorometa cresol	U	3.00	ug/L	3.00	10.0
	<i>4-Chloro-3-methylphenol</i>					
106-47-8	4-Chloroaniline	U	3.30	ug/L	3.30	10.0
7005-72-3	4-Chlorophenylphenylether	U	3.00	ug/L	3.00	10.0
100-02-7	4-Nitrophenol	U	3.00	ug/L	3.00	10.0
83-32-9	Acenaphthene	U	0.300	ug/L	0.300	1.00
208-96-8	Acenaphthylene	U	0.300	ug/L	0.300	1.00
62-53-3	Aniline	U	4.20	ug/L	4.20	10.0
120-12-7	Anthracene	U	0.300	ug/L	0.300	1.00
1912-24-9	Atrazine	U	3.00	ug/L	3.00	10.0
92-87-5	Benzidine	U	3.90	ug/L	3.90	10.0
56-55-3	Benzo(a)anthracene	U	0.300	ug/L	0.300	1.00
50-32-8	Benzo(a)pyrene	U	0.300	ug/L	0.300	1.00
205-99-2	Benzo(b)fluoranthene	U	0.300	ug/L	0.300	1.00
191-24-2	Benzo(ghi)perylene	U	0.300	ug/L	0.300	1.00

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SDG Number: 2018-1811
Lab Sample ID: 1203978892
Client Sample: QC for batch 1742247
Client ID: MB for batch 1742247
Batch ID: 1742249
Run Date: 02/27/2018 18:06
Prep Date: 02/27/2018 08:00
Data File: s022718.s\3b2713.D

Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 1000 mL
Column: DB-5ms

Matrix: WATER
Project: QC
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
207-08-9	Benzo(k)fluoranthene	U	0.300	ug/L	0.300	1.00
65-85-0	Benzoic acid	U	6.00	ug/L	6.00	20.0
100-51-6	Benzyl alcohol	U	3.00	ug/L	3.00	10.0
85-68-7	Butylbenzylphthalate	U	3.00	ug/L	3.00	10.0
218-01-9	Chrysene	U	0.300	ug/L	0.300	1.00
84-74-2	Di-n-butylphthalate	U	3.00	ug/L	3.00	10.0
117-84-0	Di-n-octylphthalate	U	3.00	ug/L	3.00	10.0
53-70-3	Dibenzo(a,h)anthracene	U	0.300	ug/L	0.300	1.00
132-64-9	Dibenzofuran	U	3.00	ug/L	3.00	10.0
84-66-2	Diethylphthalate	U	3.00	ug/L	3.00	10.0
131-11-3	Dimethylphthalate	U	3.00	ug/L	3.00	10.0
88-85-7	Dinoseb	U	3.00	ug/L	3.00	10.0
122-39-4	Diphenylamine	U	3.00	ug/L	3.00	10.0
206-44-0	Fluoranthene	U	0.300	ug/L	0.300	1.00
86-73-7	Fluorene	U	0.300	ug/L	0.300	1.00
118-74-1	Hexachlorobenzene	U	3.00	ug/L	3.00	10.0
87-68-3	Hexachlorobutadiene	U	3.00	ug/L	3.00	10.0
77-47-4	Hexachlorocyclopentadiene	U	3.00	ug/L	3.00	10.0
67-72-1	Hexachloroethane	U	3.00	ug/L	3.00	10.0
193-39-5	Indeno(1,2,3-cd)pyrene	U	0.300	ug/L	0.300	1.00
78-59-1	Isophorone	U	3.50	ug/L	3.50	10.0
62-75-9	N-Methyl-N-nitrosomethylamine	U	3.00	ug/L	3.00	10.0
924-16-3	N-Nitrosodi-n-butylamine	U	3.00	ug/L	3.00	10.0
55-18-5	N-Nitrosodiethylamine	U	3.00	ug/L	3.00	10.0
621-64-7	N-Nitrosodi-n-propylamine	U	3.00	ug/L	3.00	10.0
	<i>N-Nitrosodipropylamine</i>					
930-55-2	N-Nitrosopyrrolidine	U	3.00	ug/L	3.00	10.0
91-20-3	Naphthalene	U	0.300	ug/L	0.300	1.00
98-95-3	Nitrobenzene	U	3.00	ug/L	3.00	10.0
608-93-5	Pentachlorobenzene	U	3.00	ug/L	3.00	10.0
87-86-5	Pentachlorophenol	U	3.00	ug/L	3.00	10.0
85-01-8	Phenanthrene	U	0.300	ug/L	0.300	1.00
108-95-2	Phenol	U	3.00	ug/L	3.00	10.0
129-00-0	Pyrene	U	0.300	ug/L	0.300	1.00
110-86-1	Pyridine	U	3.00	ug/L	3.00	10.0
108-60-1	bis(2-Chloro-1-methylethyl)ether	U	3.00	ug/L	3.00	10.0
111-91-1	bis(2-Chloroethoxy)methane	U	3.00	ug/L	3.00	10.0
111-44-4	bis(2-Chloroethyl) ether	U	3.00	ug/L	3.00	10.0
117-81-7	bis(2-Ethylhexyl)phthalate	U	3.00	ug/L	3.00	1.00

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SDG Number: 2018-1811	Matrix: WATER
Lab Sample ID: 1203978892	
Client Sample: QC for batch 1742247	Client: ARSL004
Client ID: MB for batch 1742247	Method: SW846 3510C/8270D
Batch ID: 1742249	Inst: MSD3.I
Run Date: 02/27/2018 18:06	Analyst: JLD1
Prep Date: 02/27/2018 08:00	Aliquot: 1000 mL
Data File: s022718.s\s3b2713.D	Column: DB-5ms
	Project: QC
	SOP Ref: GL-OA-E-009
	Dilution: 1
	Inj. Vol: 1 uL
	Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
65794-96-9	m,p-Cresols	U	3.70	ug/L	3.70	10.0
99-09-2	3-Nitroaniline	U	3.00	ug/L	3.00	10.0
	<i>m-Nitroaniline</i>					
95-48-7	o-Cresol	U	3.00	ug/L	3.00	10.0
88-74-4	2-Nitroaniline	U	3.00	ug/L	3.00	10.0
	<i>o-Nitroaniline</i>					
100-01-6	4-Nitroaniline	U	3.00	ug/L	3.00	10.0
	<i>p-Nitroaniline</i>					

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
2,4,6-Tribromophenol	41.4	100	ug/L	41	(32%-124%)
2-Fluorobiphenyl	29.4	50.0	ug/L	59	(32%-112%)
2-Fluorophenol	34.9	100	ug/L	35	(15%-88%)
Nitrobenzene-d5	29.5	50.0	ug/L	59	(36%-115%)
Phenol-d5	21.4	100	ug/L	21	(15%-91%)
p-Terphenyl-d14	30.3	50.0	ug/L	61	(36%-121%)

Tentatively Identified Compound Summary

CAS No.	Tentatively Identified Compound (TIC)	RT	Estimated	Units	Fit	Qual
No Tentatively Identified Compounds Found				ug/L		

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SDG Number: 2018-1811
Lab Sample ID: 1203978893
Client Sample: QC for batch 1742247
Client ID: LCS for batch 1742247
Batch ID: 1742249
Run Date: 02/27/2018 18:36
Prep Date: 02/27/2018 08:00
Data File: s022718.s\3b2714.D

Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 1000 mL
Column: DB-5ms

Matrix: WATER
Project: QC
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
95-94-3	1,2,4,5-Tetrachlorobenzene		37.0	ug/L	3.00	10.0
120-82-1	1,2,4-Trichlorobenzene		34.2	ug/L	3.00	10.0
95-50-1	1,2-Dichlorobenzene		34.8	ug/L	3.00	10.0
122-66-7	Azobenzene		38.0	ug/L	3.00	10.0
	<i>1,2-Diphenylhydrazine</i>					
541-73-1	1,3-Dichlorobenzene		33.9	ug/L	3.00	10.0
106-46-7	1,4-Dichlorobenzene		34.1	ug/L	3.00	10.0
123-91-1	1,4-Dioxane		27.6	ug/L	3.00	10.0
90-12-0	1-Methylnaphthalene		33.8	ug/L	0.300	1.00
58-90-2	2,3,4,6-Tetrachlorophenol		32.0	ug/L	3.00	10.0
95-95-4	2,4,5-Trichlorophenol		33.7	ug/L	3.00	10.0
88-06-2	2,4,6-Trichlorophenol		35.0	ug/L	3.00	10.0
120-83-2	2,4-Dichlorophenol		37.1	ug/L	3.00	10.0
105-67-9	2,4-Dimethylphenol		32.4	ug/L	3.00	10.0
51-28-5	2,4-Dinitrophenol		34.3	ug/L	5.00	20.0
121-14-2	2,4-Dinitrotoluene		33.2	ug/L	3.00	10.0
606-20-2	2,6-Dinitrotoluene		34.6	ug/L	3.00	10.0
91-58-7	2-Chloronaphthalene		34.1	ug/L	0.410	1.00
95-57-8	2-Chlorophenol		37.3	ug/L	3.00	10.0
534-52-1	2-Methyl-4,6-dinitrophenol		34.6	ug/L	3.00	10.0
91-57-6	2-Methylnaphthalene		32.6	ug/L	0.300	1.00
88-75-5	2-Nitrophenol		37.0	ug/L	3.00	10.0
91-94-1	3,3'-Dichlorobenzidine		44.1	ug/L	3.00	10.0
101-55-3	4-Bromophenylphenylether		38.1	ug/L	3.00	10.0
59-50-7	Parachlorometa cresol		36.1	ug/L	3.00	10.0
	<i>4-Chloro-3-methylphenol</i>					
106-47-8	4-Chloroaniline		49.7	ug/L	3.30	10.0
7005-72-3	4-Chlorophenylphenylether		38.6	ug/L	3.00	10.0
100-02-7	4-Nitrophenol		17.6	ug/L	3.00	10.0
83-32-9	Acenaphthene		38.4	ug/L	0.300	1.00
208-96-8	Acenaphthylene		35.4	ug/L	0.300	1.00
62-53-3	Aniline		39.9	ug/L	4.20	10.0
120-12-7	Anthracene		37.1	ug/L	0.300	1.00
1912-24-9	Atrazine		32.4	ug/L	3.00	10.0
92-87-5	Benzidine		88.9	ug/L	3.90	10.0
56-55-3	Benzo(a)anthracene		36.2	ug/L	0.300	1.00
50-32-8	Benzo(a)pyrene		37.0	ug/L	0.300	1.00
205-99-2	Benzo(b)fluoranthene		34.3	ug/L	0.300	1.00
191-24-2	Benzo(ghi)perylene		44.4	ug/L	0.300	1.00

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SDG Number: 2018-1811
Lab Sample ID: 1203978893
Client Sample: QC for batch 1742247
Client ID: LCS for batch 1742247
Batch ID: 1742249
Run Date: 02/27/2018 18:36
Prep Date: 02/27/2018 08:00
Data File: s022718.s\3b2714.D

Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 1000 mL
Column: DB-5ms

Matrix: WATER
Project: QC
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
207-08-9	Benzo(k)fluoranthene		35.6	ug/L	0.300	1.00
65-85-0	Benzoic acid		40.5	ug/L	6.00	20.0
100-51-6	Benzyl alcohol		35.2	ug/L	3.00	10.0
85-68-7	Butylbenzylphthalate		35.4	ug/L	3.00	10.0
218-01-9	Chrysene		36.7	ug/L	0.300	1.00
84-74-2	Di-n-butylphthalate		42.7	ug/L	3.00	10.0
117-84-0	Di-n-octylphthalate		39.1	ug/L	3.00	10.0
53-70-3	Dibenzo(a,h)anthracene		45.3	ug/L	0.300	1.00
132-64-9	Dibenzofuran		36.7	ug/L	3.00	10.0
84-66-2	Diethylphthalate		38.9	ug/L	3.00	10.0
131-11-3	Dimethylphthalate		38.4	ug/L	3.00	10.0
88-85-7	Dinoseb	U	3.00	ug/L	3.00	10.0
122-39-4	Diphenylamine		37.4	ug/L	3.00	10.0
206-44-0	Fluoranthene		41.3	ug/L	0.300	1.00
86-73-7	Fluorene		36.7	ug/L	0.300	1.00
118-74-1	Hexachlorobenzene		37.3	ug/L	3.00	10.0
87-68-3	Hexachlorobutadiene		34.3	ug/L	3.00	10.0
77-47-4	Hexachlorocyclopentadiene		27.8	ug/L	3.00	10.0
67-72-1	Hexachloroethane		33.3	ug/L	3.00	10.0
193-39-5	Indeno(1,2,3-cd)pyrene		43.8	ug/L	0.300	1.00
78-59-1	Isophorone		35.4	ug/L	3.50	10.0
62-75-9	N-Methyl-N-nitrosomethylamine		27.5	ug/L	3.00	10.0
924-16-3	N-Nitrosodi-n-butylamine	U	3.00	ug/L	3.00	10.0
55-18-5	N-Nitrosodiethylamine	U	3.00	ug/L	3.00	10.0
621-64-7	N-Nitrosodi-n-propylamine		43.6	ug/L	3.00	10.0
	<i>N-Nitrosodipropylamine</i>					
930-55-2	N-Nitrosopyrrolidine		40.9	ug/L	3.00	10.0
91-20-3	Naphthalene		34.2	ug/L	0.300	1.00
98-95-3	Nitrobenzene		37.0	ug/L	3.00	10.0
608-93-5	Pentachlorobenzene	U	3.00	ug/L	3.00	10.0
87-86-5	Pentachlorophenol		39.4	ug/L	3.00	10.0
85-01-8	Phenanthrene		36.9	ug/L	0.300	1.00
108-95-2	Phenol		21.4	ug/L	3.00	10.0
129-00-0	Pyrene		28.2	ug/L	0.300	1.00
110-86-1	Pyridine		31.4	ug/L	3.00	10.0
108-60-1	bis(2-Chloro-1-methylethyl)ether		41.7	ug/L	3.00	10.0
111-91-1	bis(2-Chloroethoxy)methane		37.7	ug/L	3.00	10.0
111-44-4	bis(2-Chloroethyl) ether		39.6	ug/L	3.00	10.0
117-81-7	bis(2-Ethylhexyl)phthalate		40.0	ug/L	3.00	1.00

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SDG Number: 2018-1811	Matrix: WATER
Lab Sample ID: 1203978893	
Client Sample: QC for batch 1742247	Client: ARSL004
Client ID: LCS for batch 1742247	Method: SW846 3510C/8270D
Batch ID: 1742249	Inst: MSD3.I
Run Date: 02/27/2018 18:36	Analyst: JLD1
Prep Date: 02/27/2018 08:00	Aliquot: 1000 mL
Data File: s022718.s\s3b2714.D	Column: DB-5ms
	Project: QC
	SOP Ref: GL-OA-E-009
	Dilution: 1
	Inj. Vol: 1 uL
	Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
65794-96-9	m,p-Cresols		37.4	ug/L	3.70	10.0
99-09-2	3-Nitroaniline		42.6	ug/L	3.00	10.0
	<i>m-Nitroaniline</i>					
95-48-7	o-Cresol		34.1	ug/L	3.00	10.0
88-74-4	2-Nitroaniline		37.0	ug/L	3.00	10.0
	<i>o-Nitroaniline</i>					
100-01-6	4-Nitroaniline		35.9	ug/L	3.00	10.0
	<i>p-Nitroaniline</i>					

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
2,4,6-Tribromophenol	75.3	100	ug/L	75	(32%-124%)
2-Fluorobiphenyl	34.0	50.0	ug/L	68	(32%-112%)
2-Fluorophenol	55.4	100	ug/L	55	(15%-88%)
Nitrobenzene-d5	37.7	50.0	ug/L	75	(36%-115%)
Phenol-d5	43.9	100	ug/L	44	(15%-91%)
p-Terphenyl-d14	32.5	50.0	ug/L	65	(36%-121%)

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SDG Number: 2018-1811
Lab Sample ID: 1203978894
Client Sample: QC for batch 1742247
Client ID: CAWA-18-56MS
Batch ID: 1742249
Run Date: 02/28/2018 00:30
Prep Date: 02/27/2018 08:00
Data File: s022718.s\sb2726.D

Date Collected: 02/21/2018 10:44
Date Received: 02/23/2018 09:25
Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 435 mL
Column: DB-5ms

Matrix: W
Project: QC
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
95-94-3	1,2,4,5-Tetrachlorobenzene		79.2	ug/L	6.90	23.0
120-82-1	1,2,4-Trichlorobenzene		65.1	ug/L	6.90	23.0
95-50-1	1,2-Dichlorobenzene		64.4	ug/L	6.90	23.0
122-66-7	Azobenzene		73.0	ug/L	6.90	23.0
	<i>1,2-Diphenylhydrazine</i>					
541-73-1	1,3-Dichlorobenzene		63.7	ug/L	6.90	23.0
106-46-7	1,4-Dichlorobenzene		62.7	ug/L	6.90	23.0
123-91-1	1,4-Dioxane		64.7	ug/L	6.90	23.0
90-12-0	1-Methylnaphthalene		63.8	ug/L	0.690	2.30
58-90-2	2,3,4,6-Tetrachlorophenol		60.7	ug/L	6.90	23.0
95-95-4	2,4,5-Trichlorophenol		66.7	ug/L	6.90	23.0
88-06-2	2,4,6-Trichlorophenol		71.1	ug/L	6.90	23.0
120-83-2	2,4-Dichlorophenol		69.5	ug/L	6.90	23.0
105-67-9	2,4-Dimethylphenol		59.1	ug/L	6.90	23.0
51-28-5	2,4-Dinitrophenol		72.9	ug/L	11.5	46.0
121-14-2	2,4-Dinitrotoluene		61.2	ug/L	6.90	23.0
606-20-2	2,6-Dinitrotoluene		63.3	ug/L	6.90	23.0
91-58-7	2-Chloronaphthalene		71.1	ug/L	0.943	2.30
95-57-8	2-Chlorophenol		72.1	ug/L	6.90	23.0
534-52-1	2-Methyl-4,6-dinitrophenol		69.3	ug/L	6.90	23.0
91-57-6	2-Methylnaphthalene		61.2	ug/L	0.690	2.30
88-75-5	2-Nitrophenol		70.3	ug/L	6.90	23.0
91-94-1	3,3'-Dichlorobenzidine		87.2	ug/L	6.90	23.0
101-55-3	4-Bromophenylphenylether		70.2	ug/L	6.90	23.0
59-50-7	Parachlorometa cresol		64.9	ug/L	6.90	23.0
	<i>4-Chloro-3-methylphenol</i>					
106-47-8	4-Chloroaniline		91.8	ug/L	7.59	23.0
7005-72-3	4-Chlorophenylphenylether		70.8	ug/L	6.90	23.0
100-02-7	4-Nitrophenol		46.5	ug/L	6.90	23.0
83-32-9	Acenaphthene		75.4	ug/L	0.690	2.30
208-96-8	Acenaphthylene		70.9	ug/L	0.690	2.30
62-53-3	Aniline		78.6	ug/L	9.66	23.0
120-12-7	Anthracene		69.6	ug/L	0.690	2.30
1912-24-9	Atrazine		76.1	ug/L	6.90	23.0
92-87-5	Benzidine		222	ug/L	8.97	23.0
56-55-3	Benzo(a)anthracene		69.0	ug/L	0.690	2.30
50-32-8	Benzo(a)pyrene		69.5	ug/L	0.690	2.30
205-99-2	Benzo(b)fluoranthene		64.8	ug/L	0.690	2.30
191-24-2	Benzo(ghi)perylene		77.8	ug/L	0.690	2.30

**Semi-Volatile
Certificate of Analysis
Sample Summary**

SDG Number: 2018-1811
Lab Sample ID: 1203978894
Client Sample: QC for batch 1742247
Client ID: CAWA-18-56MS
Batch ID: 1742249
Run Date: 02/28/2018 00:30
Prep Date: 02/27/2018 08:00
Data File: s022718.s\s3b2726.D

Date Collected: 02/21/2018 10:44
Date Received: 02/23/2018 09:25
Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 435 mL
Column: DB-5ms

Matrix: W
Project: QC
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
207-08-9	Benzo(k)fluoranthene		67.3	ug/L	0.690	2.30
65-85-0	Benzoic acid		106	ug/L	13.8	46.0
100-51-6	Benzyl alcohol		70.7	ug/L	6.90	23.0
85-68-7	Butylbenzylphthalate		69.9	ug/L	6.90	23.0
218-01-9	Chrysene		70.6	ug/L	0.690	2.30
84-74-2	Di-n-butylphthalate		83.9	ug/L	6.90	23.0
117-84-0	Di-n-octylphthalate		79.9	ug/L	6.90	23.0
53-70-3	Dibenzo(a,h)anthracene		78.9	ug/L	0.690	2.30
132-64-9	Dibenzofuran		70.9	ug/L	6.90	23.0
84-66-2	Diethylphthalate		70.3	ug/L	6.90	23.0
131-11-3	Dimethylphthalate		70.5	ug/L	6.90	23.0
88-85-7	Dinoseb	U	6.90	ug/L	6.90	23.0
122-39-4	Diphenylamine		70.7	ug/L	6.90	23.0
206-44-0	Fluoranthene		79.8	ug/L	0.690	2.30
86-73-7	Fluorene		67.8	ug/L	0.690	2.30
118-74-1	Hexachlorobenzene		70.2	ug/L	6.90	23.0
87-68-3	Hexachlorobutadiene		63.4	ug/L	6.90	23.0
77-47-4	Hexachlorocyclopentadiene		61.4	ug/L	6.90	23.0
67-72-1	Hexachloroethane		60.4	ug/L	6.90	23.0
193-39-5	Indeno(1,2,3-cd)pyrene		77.1	ug/L	0.690	2.30
78-59-1	Isophorone		65.1	ug/L	8.05	23.0
62-75-9	N-Methyl-N-nitrosomethylamine		65.7	ug/L	6.90	23.0
924-16-3	N-Nitrosodi-n-butylamine	U	6.90	ug/L	6.90	23.0
55-18-5	N-Nitrosodiethylamine	U	6.90	ug/L	6.90	23.0
621-64-7	N-Nitrosodi-n-propylamine		77.6	ug/L	6.90	23.0
	<i>N-Nitrosodipropylamine</i>					
930-55-2	N-Nitrosopyrrolidine		75.7	ug/L	6.90	23.0
91-20-3	Naphthalene		67.3	ug/L	0.690	2.30
98-95-3	Nitrobenzene		71.8	ug/L	6.90	23.0
608-93-5	Pentachlorobenzene	U	6.90	ug/L	6.90	23.0
87-86-5	Pentachlorophenol		79.0	ug/L	6.90	23.0
85-01-8	Phenanthrene		70.0	ug/L	0.690	2.30
108-95-2	Phenol		50.3	ug/L	6.90	23.0
129-00-0	Pyrene		56.2	ug/L	0.690	2.30
110-86-1	Pyridine		71.9	ug/L	6.90	23.0
108-60-1	bis(2-Chloro-1-methylethyl)ether		78.0	ug/L	6.90	23.0
111-91-1	bis(2-Chloroethoxy)methane		69.3	ug/L	6.90	23.0
111-44-4	bis(2-Chloroethyl) ether		74.6	ug/L	6.90	23.0
117-81-7	bis(2-Ethylhexyl)phthalate		78.3	ug/L	6.90	2.30

**Semi-Volatile
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Sample Summary**

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SDG Number: 2018-1811	Date Collected: 02/21/2018 10:44	Matrix: W
Lab Sample ID: 1203978894	Date Received: 02/23/2018 09:25	
Client Sample: QC for batch 1742247	Client: ARSL004	Project: QC
Client ID: CAWA-18-56MS	Method: SW846 3510C/8270D	SOP Ref: GL-OA-E-009
Batch ID: 1742249	Inst: MSD3.I	Dilution: 1
Run Date: 02/28/2018 00:30	Analyst: JLD1	Inj. Vol: 1 uL
Prep Date: 02/27/2018 08:00	Aliquot: 435 mL	Final Volume: 1 mL
Data File: s022718.s\s3b2726.D	Column: DB-5ms	

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
65794-96-9	m,p-Cresols		76.6	ug/L	8.51	23.0
99-09-2	3-Nitroaniline		81.1	ug/L	6.90	23.0
	<i>m-Nitroaniline</i>					
95-48-7	o-Cresol		68.6	ug/L	6.90	23.0
88-74-4	2-Nitroaniline		72.4	ug/L	6.90	23.0
	<i>o-Nitroaniline</i>					
100-01-6	4-Nitroaniline		71.4	ug/L	6.90	23.0
	<i>p-Nitroaniline</i>					

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
2,4,6-Tribromophenol	138	230	ug/L	60	(32%-124%)
2-Fluorobiphenyl	74.5	115	ug/L	65	(32%-112%)
2-Fluorophenol	131	230	ug/L	57	(15%-88%)
Nitrobenzene-d5	73.9	115	ug/L	64	(36%-115%)
Phenol-d5	102	230	ug/L	45	(15%-91%)
p-Terphenyl-d14	66.0	115	ug/L	57	(36%-121%)

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811
Lab Sample ID: 1203978895
Client Sample: QC for batch 1742247
Client ID: CAWA-18-56MSD
Batch ID: 1742249
Run Date: 02/28/2018 01:00
Prep Date: 02/27/2018 08:00
Data File: s022718.s\3b2727.D

Date Collected: 02/21/2018 10:44
Date Received: 02/23/2018 09:25
Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 435 mL
Column: DB-5ms

Matrix: W
Project: QC
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
95-94-3	1,2,4,5-Tetrachlorobenzene		74.8	ug/L	6.90	23.0
120-82-1	1,2,4-Trichlorobenzene		67.3	ug/L	6.90	23.0
95-50-1	1,2-Dichlorobenzene		65.8	ug/L	6.90	23.0
122-66-7	Azobenzene		76.5	ug/L	6.90	23.0
	<i>1,2-Diphenylhydrazine</i>					
541-73-1	1,3-Dichlorobenzene		63.4	ug/L	6.90	23.0
106-46-7	1,4-Dichlorobenzene		63.9	ug/L	6.90	23.0
123-91-1	1,4-Dioxane		60.6	ug/L	6.90	23.0
90-12-0	1-Methylnaphthalene		70.4	ug/L	0.690	2.30
58-90-2	2,3,4,6-Tetrachlorophenol		64.0	ug/L	6.90	23.0
95-95-4	2,4,5-Trichlorophenol		69.3	ug/L	6.90	23.0
88-06-2	2,4,6-Trichlorophenol		71.1	ug/L	6.90	23.0
120-83-2	2,4-Dichlorophenol		75.2	ug/L	6.90	23.0
105-67-9	2,4-Dimethylphenol		60.4	ug/L	6.90	23.0
51-28-5	2,4-Dinitrophenol		79.4	ug/L	11.5	46.0
121-14-2	2,4-Dinitrotoluene		64.2	ug/L	6.90	23.0
606-20-2	2,6-Dinitrotoluene		66.8	ug/L	6.90	23.0
91-58-7	2-Chloronaphthalene		69.7	ug/L	0.943	2.30
95-57-8	2-Chlorophenol		74.3	ug/L	6.90	23.0
534-52-1	2-Methyl-4,6-dinitrophenol		73.5	ug/L	6.90	23.0
91-57-6	2-Methylnaphthalene		68.3	ug/L	0.690	2.30
88-75-5	2-Nitrophenol		75.4	ug/L	6.90	23.0
91-94-1	3,3'-Dichlorobenzidine		91.1	ug/L	6.90	23.0
101-55-3	4-Bromophenylphenylether		72.7	ug/L	6.90	23.0
59-50-7	Parachlorometa cresol		76.0	ug/L	6.90	23.0
	<i>4-Chloro-3-methylphenol</i>					
106-47-8	4-Chloroaniline		90.5	ug/L	7.59	23.0
7005-72-3	4-Chlorophenylphenylether		74.8	ug/L	6.90	23.0
100-02-7	4-Nitrophenol		45.5	ug/L	6.90	23.0
83-32-9	Acenaphthene		75.5	ug/L	0.690	2.30
208-96-8	Acenaphthylene		71.9	ug/L	0.690	2.30
62-53-3	Aniline		73.9	ug/L	9.66	23.0
120-12-7	Anthracene		72.1	ug/L	0.690	2.30
1912-24-9	Atrazine		76.7	ug/L	6.90	23.0
92-87-5	Benzidine		202	ug/L	8.97	23.0
56-55-3	Benzo(a)anthracene		71.9	ug/L	0.690	2.30
50-32-8	Benzo(a)pyrene		73.7	ug/L	0.690	2.30
205-99-2	Benzo(b)fluoranthene		68.9	ug/L	0.690	2.30
191-24-2	Benzo(ghi)perylene		74.5	ug/L	0.690	2.30

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811
Lab Sample ID: 1203978895
Client Sample: QC for batch 1742247
Client ID: CAWA-18-56MSD
Batch ID: 1742249
Run Date: 02/28/2018 01:00
Prep Date: 02/27/2018 08:00
Data File: s022718.s\s3b2727.D

Date Collected: 02/21/2018 10:44
Date Received: 02/23/2018 09:25
Client: ARSL004
Method: SW846 3510C/8270D
Inst: MSD3.I
Analyst: JLD1
Aliquot: 435 mL
Column: DB-5ms

Matrix: W
Project: QC
SOP Ref: GL-OA-E-009
Dilution: 1
Inj. Vol: 1 uL
Final Volume: 1 mL

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
207-08-9	Benzo(k)fluoranthene		70.4	ug/L	0.690	2.30
65-85-0	Benzoic acid		116	ug/L	13.8	46.0
100-51-6	Benzyl alcohol		75.2	ug/L	6.90	23.0
85-68-7	Butylbenzylphthalate		71.8	ug/L	6.90	23.0
218-01-9	Chrysene		73.6	ug/L	0.690	2.30
84-74-2	Di-n-butylphthalate		84.8	ug/L	6.90	23.0
117-84-0	Di-n-octylphthalate		83.7	ug/L	6.90	23.0
53-70-3	Dibenzo(a,h)anthracene		77.7	ug/L	0.690	2.30
132-64-9	Dibenzofuran		73.4	ug/L	6.90	23.0
84-66-2	Diethylphthalate		72.6	ug/L	6.90	23.0
131-11-3	Dimethylphthalate		73.1	ug/L	6.90	23.0
88-85-7	Dinoseb	U	6.90	ug/L	6.90	23.0
122-39-4	Diphenylamine		74.3	ug/L	6.90	23.0
206-44-0	Fluoranthene		81.0	ug/L	0.690	2.30
86-73-7	Fluorene		71.8	ug/L	0.690	2.30
118-74-1	Hexachlorobenzene		72.2	ug/L	6.90	23.0
87-68-3	Hexachlorobutadiene		63.3	ug/L	6.90	23.0
77-47-4	Hexachlorocyclopentadiene		54.7	ug/L	6.90	23.0
67-72-1	Hexachloroethane		60.8	ug/L	6.90	23.0
193-39-5	Indeno(1,2,3-cd)pyrene		76.3	ug/L	0.690	2.30
78-59-1	Isophorone		69.9	ug/L	8.05	23.0
62-75-9	N-Methyl-N-nitrosomethylamine		62.8	ug/L	6.90	23.0
924-16-3	N-Nitrosodi-n-butylamine	U	6.90	ug/L	6.90	23.0
55-18-5	N-Nitrosodiethylamine	U	6.90	ug/L	6.90	23.0
621-64-7	N-Nitrosodi-n-propylamine		85.4	ug/L	6.90	23.0
	<i>N-Nitrosodipropylamine</i>					
930-55-2	N-Nitrosopyrrolidine		84.6	ug/L	6.90	23.0
91-20-3	Naphthalene		70.7	ug/L	0.690	2.30
98-95-3	Nitrobenzene		75.7	ug/L	6.90	23.0
608-93-5	Pentachlorobenzene	U	6.90	ug/L	6.90	23.0
87-86-5	Pentachlorophenol		85.0	ug/L	6.90	23.0
85-01-8	Phenanthrene		72.9	ug/L	0.690	2.30
108-95-2	Phenol		48.6	ug/L	6.90	23.0
129-00-0	Pyrene		57.7	ug/L	0.690	2.30
110-86-1	Pyridine		64.9	ug/L	6.90	23.0
108-60-1	bis(2-Chloro-1-methylethyl)ether		81.8	ug/L	6.90	23.0
111-91-1	bis(2-Chloroethoxy)methane		74.3	ug/L	6.90	23.0
111-44-4	bis(2-Chloroethyl) ether		77.8	ug/L	6.90	23.0
117-81-7	bis(2-Ethylhexyl)phthalate		81.1	ug/L	6.90	2.30

**Semi-Volatile
Certificate of Analysis
Sample Summary**

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SDG Number: 2018-1811	Date Collected: 02/21/2018 10:44	Matrix: W
Lab Sample ID: 1203978895	Date Received: 02/23/2018 09:25	
Client Sample: QC for batch 1742247	Client: ARSL004	Project: QC
Client ID: CAWA-18-56MSD	Method: SW846 3510C/8270D	SOP Ref: GL-OA-E-009
Batch ID: 1742249	Inst: MSD3.I	Dilution: 1
Run Date: 02/28/2018 01:00	Analyst: JLD1	Inj. Vol: 1 uL
Prep Date: 02/27/2018 08:00	Aliquot: 435 mL	Final Volume: 1 mL
Data File: s022718.s\s3b2727.D	Column: DB-5ms	

CAS No.	Parmname	Qualifier	Result	Units	MDL/LOD	PQL/LOQ
65794-96-9	m,p-Cresols		83.0	ug/L	8.51	23.0
99-09-2	3-Nitroaniline		82.4	ug/L	6.90	23.0
	<i>m</i> -Nitroaniline					
95-48-7	o-Cresol		72.8	ug/L	6.90	23.0
88-74-4	2-Nitroaniline		76.3	ug/L	6.90	23.0
	<i>o</i> -Nitroaniline					
100-01-6	4-Nitroaniline		75.3	ug/L	6.90	23.0
	<i>p</i> -Nitroaniline					

Surrogate/Tracer recovery	Result	Nominal		Recovery%	Acceptable Limits
2,4,6-Tribromophenol	148	230	ug/L	64	(32%-124%)
2-Fluorobiphenyl	70.8	115	ug/L	62	(32%-112%)
2-Fluorophenol	126	230	ug/L	55	(15%-88%)
Nitrobenzene-d5	76.7	115	ug/L	67	(36%-115%)
Phenol-d5	98.6	230	ug/L	43	(15%-91%)
p-Terphenyl-d14	66.9	115	ug/L	58	(36%-121%)

Perchlorates by LCMSMS Analysis

Case Narrative

**Perchlorates by LCMSMS
Technical Case Narrative
ARS International, LLC (ARSL)
SDG #: 2018-1811
Work Order #: 444566**

Method/Analysis Information

Procedure: **Definitive Low Level Perchlorate Analysis Utilizing Liquid Chromatography/Mass Spectrometry/Mass Spectrometry (LC/MS/MS) by EPA Method 6850 Modified (6850M)**

Analytical Method: SW-846:6850

Prep Method: SW-846:6850

Analytical Batch Number: 1742510

Prep Batch Number: 1742509

Sample Analysis

Sample ID	Client ID
444566001	444566001 (CAPA-18-1)
444566005	444566005 (CAPA-18-10)
444566011	444566011 (CAWA-18-55)
1203979483	Interference Check Sample (ICS)
1203979479	Method Blank (MB)
1203979480	Laboratory Control Sample (LCS)
1203979481	444396001(CAWA-18-27) Matrix Spike (MS)
1203979482	444396001(CAWA-18-27) Matrix Spike Duplicate (MSD)

The samples in this SDG were analyzed on an "as received" basis.

Preparation/Analytical Method Verification

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-OA-E-067 REV# 14.

Calibration Information

Initial Calibration

All initial calibration requirements have been met for this SDG. Due to software constraints, all Initial Calibration Blanks must be designated as IPB001.

ICV Requirements

All associated initial calibration verification standard(s) (ICV) met the acceptance criteria.

CCB Requirements

All continuing calibration blanks (CCB) bracketing the analyses associated with this batch were within acceptance criteria.

CCV Requirements

All continuing calibration checks (CCV) requirements were met by all bracketing CCV standards.

Low Level Standard (CRI) Requirements

All low level calibration verification (CRI) requirements were met by all bracketing CRI standards.

Quality Control (QC) Information

Method Blank (MB) Statement

The MB analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries met the acceptance limits.

Interference Check Sample (ICS)

The ICS spike recoveries met the acceptance criteria.

QC Sample Designation

Client sample 444396001 (CAWA-18-27) was chosen for matrix spike and matrix spike duplicate analysis.

Matrix Spike (MS) Recovery Statement

One or more of the required spiking analytes were not within the acceptance limits in the matrix spike (MS) (See Below). Both the LCS and MSD met spike recoveries. The failing recoveries are attributed to vagaries in the extraction process. .

Sample	Value
1203979481 (CAWA-18-27MS)	59* (75%-125%)

MS/MSD Relative Percent Difference (RPD) Statement

The RPDs between the MS and MSD met the acceptance limits.

Internal Standard Area Acceptance

The internal standard areas were within the required acceptance criteria for all samples and QC.

Retention Time

During the analysis of Perchlorate by LC/MS/MS, retention time shifts are commonly observed. These retention time shifts, which are caused by fouling of the column by the sample matrices, are problematic when the retention time is used as one of the criterion for confirmation. To overcome this problem, a known amount of O(18) labeled Perchlorate was added to each sample as a retention time standard. The presence of Perchlorate was confirmed by the relative retention time (RRT) of the Perchlorate peak and the O(18) standard. A RRT window of 0.98 to 1.02, as required by DOD QSM 5.0, has been used. In addition to the isotopic ratio, the

presence of Perchlorate in the samples associated with this data package have been confirmed using the relative retention criteria stated above, not the absolute retention time.

Technical Information

Holding Time Specifications

All samples in this SDG in this analytical batch met the specified holding time. GEL assigns holding times based on the associated methodology, which assigns the date and time from sample collection of sample receipt. Those holding times expressed in hours are calculated in the AlphaLIMS system. Those holding times expressed as days expire at midnight on the day of expiration.

Preparation/Analytical Method Verification

All procedures were performed as stated in the SOP.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-extraction/Re-analysis

Re-extractions or re-analyses were not required in this SDG.

Miscellaneous Information

Manual Integrations

Manual integrations were not required for any data file associated with this SDG.

Method Comments

The samples in this SDG were not originally analyzed using EPA Method 314.0.

Additional Comments

The Perchlorate Isotope Ratio on the Form I may differ slightly from the ratio on the corresponding raw data due to rounding rules and/or significant figures or due to software limitations when there are manual integrations, dilutions or other factors. The ratio value of the Form I is the correct value. The retention time marker, Perchlorate-O (18), is added to all samples, instrument blanks, and standards prior to injection. It is used to verify the retention time of Perchlorate and Perchlorate-101 and to insure an accurate injection occurred. Due to various anions affecting the recovery of Perchlorate-O (18) and not Perchlorate and Perchlorate-101, the calibration curves of Perchlorate and Perchlorate-101 are internally corrected for using Perchlorate-O (18).

Perchlorate Isotope Ratio

The Perchlorate isotope ratio met acceptance criteria for all samples and QC samples. Please see the isotope ratio criteria in the Miscellaneous Section.

System Configuration

The laboratory utilizes a Waters LC 2795 liquid chromatography instrument for Perchlorate analysis. It is coupled with a Micromass Quattro Ultima Mass Spectrometer/Mass Spectrometer. It is designated as LCMSMS #2. It is fitted with an electrospray probe that is operated in the negative electrospray ionization mode for Perchlorate analysis. The laboratory may also utilize an Agilent 1100 liquid chromatography instrument for Perchlorate analysis. It is coupled with an Applied Biosystems 4000 Mass Spectrometer/Mass Spectrometer, designated as LCMSMS #3 or LCMSMS #4. It is also fitted with an electrospray probe that is operated in the negative electrospray ionization mode for Perchlorate analysis.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. An electronic signature page inserted after the case narrative will include the data validator's signature and title. The signature page also includes the data qualifiers used in the fractional package. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Chromatographic Columns

The LC-MS/MS Perchlorate analysis was performed on a Quatro Ultima LC/MS/MS.

Chromatographic separation of Perchlorate is accomplished through analysis on the following anion column:

Dionex: IonPac AG-16 2 x 50 mm.

Certification Statement

Where the analytical method has been performed under NELAP certification, the analysis has met all of the requirements of the NELAC standard unless otherwise noted in the analytical case narrative.

GEL LABORATORIES LLC

2040 Savage Road Charleston SC 29407 - (843) 556-8171 - www.gel.com

Qualifier Definition Report for

ARSL004 ARS International, LLC (ARS-LANS-MTOA6-25093-GEL)

Client SDG: 2018-1811 GEL Work Order: 444566

The Qualifiers in this report are defined as follows:

- * A quality control analyte recovery is outside of specified acceptance criteria
- ** Analyte is a surrogate compound
- J Value is estimated
- U Analyte was analyzed for, but not detected above the MDL, MDA, MDC or LOD.
- DL Indicates that sample is diluted.
- RA Indicates that sample is re-analyzed without re-extraction.
- RE Indicates that sample is re-extracted.

Review/Validation

GEL requires all analytical data to be verified by a qualified data reviewer. In addition, all CLP-like deliverables receive a third level review of the fractional data package.

The following data validator verified the information presented in this data report:

Signature: 

Name: Michael Penny

Date: 02 MAR 2018

Title: Group Leader

Sample Data Summary

Perchlorate Analysis Data Sheet

Lab Name: GEL Laboratories LLCLab Code: GELInstrument: LCMSMSMethod: SW846 6850 ModifiedMatrix: WATERExtraction Batch ID: 1742509Extraction Type: Filter/DAISample Volume/Weight: 10.0 mLConcentrated Extract Volume: 10.0

Client Sample No.

CAPA-18-1Date Received: 23-FEB-18GEL Job No (SDG): 2018-1811GEL Sample ID: 444566001Date Filtered: 27-FEB-18Injection Volume (uL): 20%Solids:

CAS No.	Analyte^	MDL	RL	Conc*	Units	Q	Dilution Factor	Date Analyzed	GEL File ID
14797-73-0	Perchlorate	.05	.2	0.992	ug/L		1	28-FEB-18 18:01	per0228033a
	Perchlorate Isotope Ratio			2.91			1	28-FEB-18 18:01	per0228033a
14797-73-0	Perchlorate-101	.05	.2	0.993	ug/L		1	28-FEB-18 18:01	per0228033a
	Perchlorate-O(18)			0.468	ug/L		1	28-FEB-18 18:01	per0228033a

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

*Concentration =

$$\text{Instrument Value} \times \frac{\text{Concentrated Extract Volume}}{\text{Aliquot}} \times \frac{1}{\% \text{Solids}}$$

Perchlorate Analysis Data Sheet

Lab Name: GEL Laboratories LLCLab Code: GELInstrument: LCMSMSMethod: SW846 6850 ModifiedMatrix: WATERExtraction Batch ID: 1742509Extraction Type: Filter/DAISample Volume/Weight: 10.0 mLConcentrated Extract Volume: 10.0

Client Sample No.

CAPA-18-10Date Received: 23-FEB-18GEL Job No (SDG): 2018-1811GEL Sample ID: 444566005Date Filtered: 27-FEB-18Injection Volume (uL): 20%Solids:

CAS No.	Analyte^	MDL	RL	Conc*	Units	Q	Dilution Factor	Date Analyzed	GEL File ID
14797-73-0	Perchlorate	.05	.2	0.968	ug/L		1	28-FEB-18 18:09	per0228034a
	Perchlorate Isotope Ratio			2.83			1	28-FEB-18 18:09	per0228034a
14797-73-0	Perchlorate-101	.05	.2	0.997	ug/L		1	28-FEB-18 18:09	per0228034a
	Perchlorate-O(18)			0.461	ug/L		1	28-FEB-18 18:09	per0228034a

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

*Concentration =

$$\text{Instrument Value} \times \frac{\text{Concentrated Extract Volume}}{\text{Aliquot}} \times \frac{1}{\% \text{Solids}}$$

Perchlorate Analysis Data Sheet

Lab Name: GEL Laboratories LLCLab Code: GELInstrument: LCMSMSMethod: SW846 6850 ModifiedMatrix: WATERExtraction Batch ID: 1742509Extraction Type: Filter/DAISample Volume/Weight: 10.0 mLConcentrated Extract Volume: 10.0

Client Sample No.

CAWA-18-55Date Received: 23-FEB-18GEL Job No (SDG): 2018-1811GEL Sample ID: 444566011Date Filtered: 27-FEB-18Injection Volume (uL): 20%Solids:

CAS No.	Analyte^	MDL	RL	Conc*	Units	Q	Dilution Factor	Date Analyzed	GEL File ID
14797-73-0	Perchlorate	.05	.2	0.254	ug/L		1	28-FEB-18 18:17	per0228035a
	Perchlorate Isotope Ratio			2.99			1	28-FEB-18 18:17	per0228035a
14797-73-0	Perchlorate-101	.05	.2	0.248	ug/L		1	28-FEB-18 18:17	per0228035a
	Perchlorate-O(18)			0.401	ug/L		1	28-FEB-18 18:17	per0228035a

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

*Concentration =

$$\text{Instrument Value} \times \frac{\text{Concentrated Extract Volume}}{\text{Aliquot}} \times \frac{1}{\% \text{Solids}}$$

Quality Control Summary

Perchlorate Laboratory Control Sample

Lab Name: General Engineering Laboratories

Lab Code: GEL

GEL Job No. (SDG): 2018-1811

Extract Batch Code: 1742509

Date Filtered: 27-FEB-18

Matrix: WATER

Sample ID: 1203979480

Analyte^	True	Found	Units	%Rec	Q	Control Limits
Perchlorate	0.200	.176	ug/L	88		85 - 115
Perchlorate Isotope Ratio		2.76				-
Perchlorate-101	0.200	.186	ug/L	93		85 - 115
Perchlorate-O(18)		.479	ug/L			-

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

Perchlorate Spike/Spike Duplicate Summary

Lab Name: General Engineering Laboratories

Lab Code: GEL

GEL Job No (SDG): 2018-1811

Extract Batch Code: 1742509

Date Extracted: 27-FEB-18

GEL MS/PS ID: 1203979481

Client ID: CAWA-18-27

GEL MSD/PSD ID: 1203979482

QC Type: MS

Compound^	Spike Added	Sample Conc	Units	MS Conc	MS Rec #	MSD Conc	MSD Rec #	RPD #	RPD Limit	Recovery Limit
Perchlorate	0.200	0.401	ug/L	0.520	59 *	.6	100	14	30	75 - 125
Perchlorate Isotope Ratio	0	3.23		2.78		3.01		8		-
Perchlorate-101	0.200	0.362	ug/L	0.546	92	.581	110	6	30	75 - 125
Perchlorate-O(18)	0	0.464	ug/L	0.467		.409		13		-

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

Quality Control Data

Perchlorate Analysis Data Sheet

Lab Name: GEL Laboratories LLCLab Code: GELInstrument: LCMSMSMethod: EPA 6850 ModifiedMatrix: WATERExtraction Batch ID: 1742509Extraction Type: Filter/DAISample Volume/Weight: 10.0 mLConcentrated Extract Volume: 10.0

Client Sample No.

MBDate Received: 27-FEB-18GEL Job No (SDG): 2018-1811GEL Sample ID: 1203979479Date Filtered: 27-FEB-18Injection Volume (uL): 20%Solids:

CAS No.	Analyte^	MDL	RL	Conc*	Units	Q	Dilution Factor	Date Analyzed	GEL File ID
14797-73-0	Perchlorate	.05	.2	0.050	ug/L	U	1	28-FEB-18 15:22	per0228013a
	Perchlorate Isotope Ratio						1	28-FEB-18 15:22	per0228013a
14797-73-0	Perchlorate-101	.05	.2	0.050	ug/L	U	1	28-FEB-18 15:22	per0228013a
	Perchlorate-O(18)			0.449	ug/L		1	28-FEB-18 15:22	per0228013a

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

*Concentration =

$$\text{Instrument Value} \times \frac{\text{Concentrated Extract Volume}}{\text{Aliquot}} \times \frac{1}{\% \text{Solids}}$$

Perchlorate Analysis Data Sheet

Lab Name: GEL Laboratories LLCLab Code: GELInstrument: LCMSMSMethod: EPA 6850 ModifiedMatrix: WATERExtraction Batch ID: 1742509Extraction Type: Filter/DAISample Volume/Weight: 10.0 mLConcentrated Extract Volume: 10.0

Client Sample No.

LCSDate Received: 27-FEB-18GEL Job No (SDG): 2018-1811GEL Sample ID: 1203979480Date Filtered: 27-FEB-18Injection Volume (uL): 20%Solids:

CAS No.	Analyte^	MDL	RL	Conc*	Units	Q	Dilution Factor	Date Analyzed	GEL File ID
14797-73-0	Perchlorate	.05	.2	0.176	ug/L	J	1	28-FEB-18 15:30	per0228014a
	Perchlorate Isotope Ratio			2.76			1	28-FEB-18 15:30	per0228014a
14797-73-0	Perchlorate-101	.05	.2	0.186	ug/L	J	1	28-FEB-18 15:30	per0228014a
	Perchlorate-O(18)			0.479	ug/L		1	28-FEB-18 15:30	per0228014a

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

*Concentration =

$$\text{Instrument Value} \times \frac{\text{Concentrated Extract Volume}}{\text{Aliquot}} \times \frac{1}{\% \text{Solids}}$$

Perchlorate Analysis Data Sheet

Lab Name: GEL Laboratories LLCLab Code: GELInstrument: LCMSMSMethod: SW846 6850 ModifiedMatrix: WATERExtraction Batch ID: 1742509Extraction Type: Filter/DAISample Volume/Weight: 10.0 mLConcentrated Extract Volume: 10.0

Client Sample No.

ICS

Date Received:

GEL Job No (SDG): 2018-1811GEL Sample ID: 1203979483Date Filtered: 27-FEB-18Injection Volume (uL): 20

%Solids:

CAS No.	Analyte^	MDL	RL	Conc*	Units	Q	Dilution Factor	Date Analyzed	GEL File ID
14797-73-0	Perchlorate	.05	.2	0.213	ug/L		1	28-FEB-18 15:38	per0228015a
	Perchlorate Isotope Ratio			2.85			1	28-FEB-18 15:38	per0228015a
14797-73-0	Perchlorate-101	.05	.2	0.217	ug/L		1	28-FEB-18 15:38	per0228015a
	Perchlorate-O(18)			0.483	ug/L		1	28-FEB-18 15:38	per0228015a

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

*Concentration =

$$\text{Instrument Value} \times \frac{\text{Concentrated Extract Volume}}{\text{Aliquot}} \times \frac{1}{\% \text{Solids}}$$

Perchlorate Analysis Data Sheet

Lab Name: GEL Laboratories LLCLab Code: GELInstrument: LCMSMSMethod: SW846 6850 ModifiedMatrix: WATERExtraction Batch ID: 1742509Extraction Type: Filter/DAISample Volume/Weight: 10.0 mLConcentrated Extract Volume: 10.0

Client Sample No.

CAWA-18-27MSDate Received: 21-FEB-18GEL Job No (SDG): 2018-1811GEL Sample ID: 1203979481Date Filtered: 27-FEB-18Injection Volume (uL): 20%Solids:

CAS No.	Analyte^	MDL	RL	Conc*	Units	Q	Dilution Factor	Date Analyzed	GEL File ID
14797-73-0	Perchlorate	.05	.2	0.520	ug/L		1	28-FEB-18 15:53	per0228017a
	Perchlorate Isotope Ratio			2.78			1	28-FEB-18 15:53	per0228017a
14797-73-0	Perchlorate-101	.05	.2	0.546	ug/L		1	28-FEB-18 15:53	per0228017a
	Perchlorate-O(18)			0.467	ug/L		1	28-FEB-18 15:53	per0228017a

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

*Concentration =

$$\text{Instrument Value} \times \frac{\text{Concentrated Extract Volume}}{\text{Aliquot}} \times \frac{1}{\% \text{Solids}}$$

Perchlorate Analysis Data Sheet

Lab Name: GEL Laboratories LLCLab Code: GELInstrument: LCMSMSMethod: SW846 6850 ModifiedMatrix: WATERExtraction Batch ID: 1742509Extraction Type: Filter/DAISample Volume/Weight: 10.0 mLConcentrated Extract Volume: 10.0

Client Sample No.

CAWA-18-27MSDDate Received: 21-FEB-18GEL Job No (SDG): 2018-1811GEL Sample ID: 1203979482Date Filtered: 27-FEB-18Injection Volume (uL): 20%Solids:

CAS No.	Analyte^	MDL	RL	Conc*	Units	Q	Dilution Factor	Date Analyzed	GEL File ID
14797-73-0	Perchlorate	.05	.2	0.600	ug/L		1	28-FEB-18 16:01	per0228018a
	Perchlorate Isotope Ratio			3.01			1	28-FEB-18 16:01	per0228018a
14797-73-0	Perchlorate-101	.05	.2	0.581	ug/L		1	28-FEB-18 16:01	per0228018a
	Perchlorate-O(18)			0.409	ug/L		1	28-FEB-18 16:01	per0228018a

^ When the analyte name is Perchlorate Isotope Ratio the concentration is a unitless value calculated from the ratio of Perchlorate peak area to Perchlorate-101 peak area. The Perchlorate-101 and isotopic ratio results are provided for qualitative purposes only. The results are used to verify the presence and quantitation of Perchlorate.

*Concentration =

$$\text{Instrument Value} \times \frac{\text{Concentrated Extract Volume}}{\text{Aliquot}} \times \frac{1}{\% \text{Solids}}$$

Explosives by LCMSMS Analysis

Case Narrative

**Explosives by LCMSMS
Technical Case Narrative
ARS International, LLC (ARSL)
SDG #: 2018-1811
Work Order #: 444566**

Method/Analysis Information

Procedure: The Processing, Extraction, and Analysis of Nitroaromatics, Nitroamines, and Nitrate Esters by SW-846 8330B

Analytical Method: SW846 3535A/8330B

Prep Method: SW846 3535A

Analytical Batch Number: 1742521

Prep Batch Number: 1742520

Sample Analysis

The following samples were analyzed using the analytical protocol as established in SW846 3535A/8330B:

Sample ID	Client ID
444566003	CAPA-18-2
444566007	CAPA-18-12
444566008	CAWA-18-54
444566013	CAWA-18-56
1203979520	Method Blank (MB)
1203979521	Laboratory Control Sample (LCS)
1203979522	444566003(CAPA-18-2) Matrix Spike (MS)
1203979523	444566003(CAPA-18-2) Matrix Spike Duplicate (MSD)

Preparation/Analytical Method Verification

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-OA-E-068 REV# 7.

Calibration Information

Initial Calibration

All initial calibration requirements for this analysis have been met for this SDG.

Calibration Verification Standard Requirements

All associated calibration verification standards (ICV and CCV) for this analysis met the acceptance criteria.

Calibration Blank Requirements

All initial and continuing calibration blanks (ICB and CCB) bracketing the analyses associated with this batch

for this analysis were within acceptance criteria. Due to software limitations, the CCBs and/or the ICBs may have a concentration for target analytes in the Found column. These values should be zero.

CRI Requirements

All low level calibration verification (CRI) requirements for this analysis were met by all bracketing CRI standards.

Quality Control (QC) Information

Method Blank (MB) Statement

The MB analyzed with this SDG for this analysis met the acceptance criteria.

Surrogate Recoveries

All the surrogate recoveries were within the established acceptance criteria in this SDG in this analytical batch for this analysis.

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries were within the established acceptance limits.

QC Sample Designation

Client sample 444566003 (CAPA-18-2) was chosen for matrix spike and matrix spike duplicate analysis.

Matrix Spike (MS) Recovery Statement

The MS spike recoveries were within the established acceptance limits for this analysis.

MS/MSD Relative Percent Difference (RPD) Statement

The RPDs between the MS and MSD met the acceptance limits for this analysis.

Internal Standard (ISTD) Acceptance

The internal standard responses were within the required acceptance criteria for all samples and QC in this SDG.

Technical Information

Holding Time Specifications

All samples in this SDG in this analytical batch met the specified holding time. GEL assigns holding times based on the associated methodology, which assigns the date and time from sample collection of sample receipt. Those holding times expressed in hours are calculated in the AlphaLIMS system. Those holding times expressed as days expire at midnight on the day of expiration. All samples in this SDG met the specified holding time.

Preparation/Analytical Method Verification

All procedures were performed as stated in the SOP.

Sample Dilutions

In accordance with GEL SOP GL-OA-056, all sample and QC extracts are diluted 1:1 v/v with LC reagent grade Water. The samples in this SDG in this analytical batch for this analysis did not require any additional dilutions.

Sample Re-extraction/Re-analysis

Re-extractions or re-analyses were not required in this SDG in this analytical batch for this analysis.

Miscellaneous Information

Manual Integrations

Manual integrations were not required for any data file associated with this SDG.

Additional Comments

Due to software limitations, all initial calibration blanks must be designated as XIB001 in order for the forms to be correct. Due to software limitations, file extensions such as DL, RE, etc. may not appear on the generated forms and/or raw data. Relative Retention Time (RRT) is used by the laboratory to establish peak identity. The RRT of each target analyte is calculated using the retention time of the corresponding internal standard. The RRT of each analyte in a sample must be within 2.0 of the analyte's calculated RRT in the ICV.

System Configuration

The laboratory utilizes an Agilent 1100 liquid chromatography instrument for either Primary or Secondary analyte analysis. It is coupled with an Applied Biosystems 4000 Mass Spectrometer/ Mass Spectrometer, designated as either LC/MS/MS #3 or LC/MS/MS #4. The laboratory also utilizes a Shimadzu Nexera XC liquid chromatography instrument for Primary and/or Secondary analyte analysis. It is coupled with an Applied Biosystems 5500 Mass Spectrometer/ Mass Spectrometer, designated as LC/MS/MS #5. All are fitted with an APCI (Atmospheric Pressure Chemical Ionization) probe that is operated in the negative ionization mode for both the Primary and Secondary analyte analysis.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. An electronic signature page inserted after the case narrative will include the data validator's signature and title. The signature page also includes the data qualifiers used in the fractional package. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Chromatographic Columns

The LC-MS/MS Explosives analysis was performed on a ABSciex 5500 Qtrap LC/MS/MS.

The detection of the Primary and Secondary Nitroaromatic and Nitramine analytes is accomplished through analysis on the following reversed phase column:

Phenomenex: Ultracarb 5u ODS (20), 250 x 4.60 mm ID.

Certification Statement

Where the analytical method has been performed under NELAP certification, the analysis has met all of the requirements of the NELAC standard unless otherwise noted in the analytical case narrative.

GEL LABORATORIES LLC

2040 Savage Road Charleston SC 29407 - (843) 556-8171 - www.gel.com

Qualifier Definition Report for

ARSL004 ARS International, LLC (ARS-LANS-MTOA6-25093-GEL)

Client SDG: 2018-1811 GEL Work Order: 444566

The Qualifiers in this report are defined as follows:

- * A quality control analyte recovery is outside of specified acceptance criteria
- ** Analyte is a surrogate compound
- J Value is estimated
- U Analyte was analyzed for, but not detected above the MDL, MDA, MDC or LOD.
- DL Indicates that sample is diluted.
- RA Indicates that sample is re-analyzed without re-extraction.
- RE Indicates that sample is re-extracted.

Review/Validation

GEL requires all analytical data to be verified by a qualified data reviewer. In addition, all CLP-like deliverables receive a third level review of the fractional data package.

The following data validator verified the information presented in this data report:

Signature: 

Name: Michael Penny

Date: 12 MAR 2018

Title: Group Leader

Sample Data Summary

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAPA-18-2

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 444566003

Sample Amount 870 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

GEL data file: EXP0305018.wiff

Date Analyzed: 05-MAR-18 20:09

Dilution Factor: 2

Concentration Units: ug/L

Cas No.	Compound	Concentration*	Q	MDL	PQL
118-96-7	2,4,6-Trinitrotoluene	.092	U	0.092	0.287
<i>118-96-7</i>	<i>2,4,6-Trinitrotoluene</i>				
121-14-2	2,4-Dinitrotoluene	.092	U	0.092	0.287
<i>121-14-2</i>	<i>2,4-Dinitrotoluene</i>				
13980-04-6	TNX	.092	U	0.092	0.287
<i>13980-04-6</i>	<i>TNX</i>				
479-45-8	Tetryl	.092	U	0.092	0.575
<i>479-45-8</i>	<i>Tetryl</i>				
5755-27-1	MXN	.092	U	0.092	0.287
<i>5755-27-1</i>	<i>MXN</i>				
606-20-2	2,6-Dinitrotoluene	.092	U	0.092	0.287
<i>606-20-2</i>	<i>2,6-Dinitrotoluene</i>				
80251-29-2	DNX	.092	U	0.092	0.287
<i>80251-29-2</i>	<i>DNX</i>				
98-95-3	Nitrobenzene	.092	U	0.092	0.287
<i>98-95-3</i>	<i>Nitrobenzene</i>				
99-08-1	m-Nitrotoluene	.092	U	0.092	0.287
<i>99-08-1</i>	<i>m-Nitrotoluene</i>				
99-35-4	1,3,5-Trinitrobenzene	.092	U	0.092	0.287
<i>99-35-4</i>	<i>1,3,5-Trinitrobenzene</i>				
99-65-0	m-Dinitrobenzene	.092	U	0.092	0.287
<i>99-65-0</i>	<i>m-Dinitrobenzene</i>				
88-72-2	o-Nitrotoluene	.0943	U	0.0943	0.287
<i>88-72-2</i>	<i>o-Nitrotoluene</i>				
78-11-5	PETN	.115	U	0.115	0.575
<i>78-11-5</i>	<i>PETN</i>				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAPA-18-2

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 444566003

Sample Amount 870 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

Cas No.	Compound	Concentration*	Q	MDL	PQL
35572-78-2	2-Amino-4,6-dinitrotoluene	.12	J	0.092	0.287
35572-78-2	2-Amino-4,6-dinitrotoluene				
99-99-0	p-Nitrotoluene	.172	U	0.172	0.575
99-99-0	p-Nitrotoluene				
19406-51-0	4-Amino-2,6-dinitrotoluene	.195	J	0.092	0.287
19406-51-0	4-Amino-2,6-dinitrotoluene				
3058-38-6	TATB	.345	U	0.345	1.15
3058-38-6	TATB				
618-87-1	3,5-Dinitroaniline	.345	U	0.345	1.15
618-87-1	3,5-Dinitroaniline				
78-30-8	tris(o-cresyl) phosphate	.345	U	0.345	1.15
78-30-8	tris(o-cresyl) phosphate				
59229-75-3	2,6-Diamino-4-nitrotoluene	.575	U	0.575	2.87
59229-75-3	2,6-Diamino-4-nitrotoluene				
6629-29-4	2,4-Diamino-6-nitrotoluene	.575	U	0.575	2.87
6629-29-4	2,4-Diamino-6-nitrotoluene				
2691-41-0	HMX	6.64		0.092	0.287
2691-41-0	HMX				
121-82-4	RDX	7.69		0.092	0.287
121-82-4	RDX				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAPA-18-12

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 444566007

Sample Amount 880 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

GEL data file: EXP0305022.wiff

Date Analyzed: 05-MAR-18 22:31

Dilution Factor: 2

Concentration Units: ug/L

Cas No.	Compound	Concentration*	Q	MDL	PQL
118-96-7	2,4,6-Trinitrotoluene	.0909	U	0.0909	0.284
<i>118-96-7</i>	<i>2,4,6-Trinitrotoluene</i>				
121-14-2	2,4-Dinitrotoluene	.0909	U	0.0909	0.284
<i>121-14-2</i>	<i>2,4-Dinitrotoluene</i>				
13980-04-6	TNX	.0909	U	0.0909	0.284
<i>13980-04-6</i>	<i>TNX</i>				
479-45-8	Tetryl	.0909	U	0.0909	0.568
<i>479-45-8</i>	<i>Tetryl</i>				
5755-27-1	MXN	.0909	U	0.0909	0.284
<i>5755-27-1</i>	<i>MXN</i>				
606-20-2	2,6-Dinitrotoluene	.0909	U	0.0909	0.284
<i>606-20-2</i>	<i>2,6-Dinitrotoluene</i>				
80251-29-2	DNX	.0909	U	0.0909	0.284
<i>80251-29-2</i>	<i>DNX</i>				
98-95-3	Nitrobenzene	.0909	U	0.0909	0.284
<i>98-95-3</i>	<i>Nitrobenzene</i>				
99-08-1	m-Nitrotoluene	.0909	U	0.0909	0.284
<i>99-08-1</i>	<i>m-Nitrotoluene</i>				
99-35-4	1,3,5-Trinitrobenzene	.0909	U	0.0909	0.284
<i>99-35-4</i>	<i>1,3,5-Trinitrobenzene</i>				
99-65-0	m-Dinitrobenzene	.0909	U	0.0909	0.284
<i>99-65-0</i>	<i>m-Dinitrobenzene</i>				
88-72-2	o-Nitrotoluene	.0932	U	0.0932	0.284
<i>88-72-2</i>	<i>o-Nitrotoluene</i>				
78-11-5	PETN	.114	U	0.114	0.568
<i>78-11-5</i>	<i>PETN</i>				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAPA-18-12

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 444566007

Sample Amount 880 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

Cas No.	Compound	Concentration*	Q	MDL	PQL
35572-78-2	2-Amino-4,6-dinitrotoluene	.122	J	0.0909	0.284
35572-78-2	2-Amino-4,6-dinitrotoluene				
99-99-0	p-Nitrotoluene	.17	U	0.170	0.568
99-99-0	p-Nitrotoluene				
19406-51-0	4-Amino-2,6-dinitrotoluene	.214	J	0.0909	0.284
19406-51-0	4-Amino-2,6-dinitrotoluene				
3058-38-6	TATB	.341	U	0.341	1.14
3058-38-6	TATB				
618-87-1	3,5-Dinitroaniline	.341	U	0.341	1.14
618-87-1	3,5-Dinitroaniline				
78-30-8	tris(o-cresyl) phosphate	.341	U	0.341	1.14
78-30-8	tris(o-cresyl) phosphate				
59229-75-3	2,6-Diamino-4-nitrotoluene	.568	U	0.568	2.84
59229-75-3	2,6-Diamino-4-nitrotoluene				
6629-29-4	2,4-Diamino-6-nitrotoluene	.568	U	0.568	2.84
6629-29-4	2,4-Diamino-6-nitrotoluene				
2691-41-0	HMX	5.75		0.0909	0.284
2691-41-0	HMX				
121-82-4	RDX	7.08		0.0909	0.284
121-82-4	RDX				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAWA-18-54

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 444566008

Sample Amount 850 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

GEL data file: EXP0305023.wiff

Date Analyzed: 05-MAR-18 23:07

Dilution Factor: 2

Concentration Units: ug/L

Cas No.	Compound	Concentration*	Q	MDL	PQL
118-96-7	2,4,6-Trinitrotoluene	.0941	U	0.0941	0.294
<i>118-96-7</i>	<i>2,4,6-Trinitrotoluene</i>				
121-14-2	2,4-Dinitrotoluene	.0941	U	0.0941	0.294
<i>121-14-2</i>	<i>2,4-Dinitrotoluene</i>				
121-82-4	RDX	.0941	U	0.0941	0.294
<i>121-82-4</i>	<i>RDX</i>				
13980-04-6	TNX	.0941	U	0.0941	0.294
<i>13980-04-6</i>	<i>TNX</i>				
19406-51-0	4-Amino-2,6-dinitrotoluene	.0941	U	0.0941	0.294
<i>19406-51-0</i>	<i>4-Amino-2,6-dinitrotoluene</i>				
2691-41-0	HMX	.0941	U	0.0941	0.294
<i>2691-41-0</i>	<i>HMX</i>				
35572-78-2	2-Amino-4,6-dinitrotoluene	.0941	U	0.0941	0.294
<i>35572-78-2</i>	<i>2-Amino-4,6-dinitrotoluene</i>				
479-45-8	Tetryl	.0941	U	0.0941	0.588
<i>479-45-8</i>	<i>Tetryl</i>				
5755-27-1	MNX	.0941	U	0.0941	0.294
<i>5755-27-1</i>	<i>MNX</i>				
606-20-2	2,6-Dinitrotoluene	.0941	U	0.0941	0.294
<i>606-20-2</i>	<i>2,6-Dinitrotoluene</i>				
80251-29-2	DNX	.0941	U	0.0941	0.294
<i>80251-29-2</i>	<i>DNX</i>				
98-95-3	Nitrobenzene	.0941	U	0.0941	0.294
<i>98-95-3</i>	<i>Nitrobenzene</i>				
99-08-1	m-Nitrotoluene	.0941	U	0.0941	0.294
<i>99-08-1</i>	<i>m-Nitrotoluene</i>				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAWA-18-54

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 444566008

Sample Amount 850 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

Cas No.	Compound	Concentration*	Q	MDL	PQL
99-35-4	1,3,5-Trinitrobenzene	.0941	U	0.0941	0.294
99-35-4	1,3,5-Trinitrobenzene				
99-65-0	m-Dinitrobenzene	.0941	U	0.0941	0.294
99-65-0	m-Dinitrobenzene				
88-72-2	o-Nitrotoluene	.0965	U	0.0965	0.294
88-72-2	o-Nitrotoluene				
78-11-5	PETN	.118	U	0.118	0.588
78-11-5	PETN				
99-99-0	p-Nitrotoluene	.176	U	0.176	0.588
99-99-0	p-Nitrotoluene				
3058-38-6	TATB	.353	U	0.353	1.18
3058-38-6	TATB				
618-87-1	3,5-Dinitroaniline	.353	U	0.353	1.18
618-87-1	3,5-Dinitroaniline				
78-30-8	tris(o-cresyl) phosphate	.353	U	0.353	1.18
78-30-8	tris(o-cresyl) phosphate				
59229-75-3	2,6-Diamino-4-nitrotoluene	.588	U	0.588	2.94
59229-75-3	2,6-Diamino-4-nitrotoluene				
6629-29-4	2,4-Diamino-6-nitrotoluene	.588	U	0.588	2.94
6629-29-4	2,4-Diamino-6-nitrotoluene				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAWA-18-56

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 444566013

Sample Amount 940 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

GEL data file: EXP0305024.wiff

Date Analyzed: 05-MAR-18 23:42

Dilution Factor: 2

Concentration Units: ug/L

Cas No.	Compound	Concentration*	Q	MDL	PQL
118-96-7	2,4,6-Trinitrotoluene	.0851	U	0.0851	0.266
<i>118-96-7</i>	<i>2,4,6-Trinitrotoluene</i>				
121-14-2	2,4-Dinitrotoluene	.0851	U	0.0851	0.266
<i>121-14-2</i>	<i>2,4-Dinitrotoluene</i>				
121-82-4	RDX	.0851	U	0.0851	0.266
<i>121-82-4</i>	<i>RDX</i>				
13980-04-6	TNX	.0851	U	0.0851	0.266
<i>13980-04-6</i>	<i>TNX</i>				
19406-51-0	4-Amino-2,6-dinitrotoluene	.0851	U	0.0851	0.266
<i>19406-51-0</i>	<i>4-Amino-2,6-dinitrotoluene</i>				
2691-41-0	HMX	.0851	U	0.0851	0.266
<i>2691-41-0</i>	<i>HMX</i>				
35572-78-2	2-Amino-4,6-dinitrotoluene	.0851	U	0.0851	0.266
<i>35572-78-2</i>	<i>2-Amino-4,6-dinitrotoluene</i>				
479-45-8	Tetryl	.0851	U	0.0851	0.532
<i>479-45-8</i>	<i>Tetryl</i>				
5755-27-1	MNX	.0851	U	0.0851	0.266
<i>5755-27-1</i>	<i>MNX</i>				
606-20-2	2,6-Dinitrotoluene	.0851	U	0.0851	0.266
<i>606-20-2</i>	<i>2,6-Dinitrotoluene</i>				
80251-29-2	DNX	.0851	U	0.0851	0.266
<i>80251-29-2</i>	<i>DNX</i>				
98-95-3	Nitrobenzene	.0851	U	0.0851	0.266
<i>98-95-3</i>	<i>Nitrobenzene</i>				
99-08-1	m-Nitrotoluene	.0851	U	0.0851	0.266
<i>99-08-1</i>	<i>m-Nitrotoluene</i>				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAWA-18-56

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 444566013

Sample Amount 940 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

Cas No.	Compound	Concentration*	Q	MDL	PQL
99-35-4	1,3,5-Trinitrobenzene	.0851	U	0.0851	0.266
99-35-4	1,3,5-Trinitrobenzene				
99-65-0	m-Dinitrobenzene	.0851	U	0.0851	0.266
99-65-0	m-Dinitrobenzene				
88-72-2	o-Nitrotoluene	.0872	U	0.0872	0.266
88-72-2	o-Nitrotoluene				
78-11-5	PETN	.106	U	0.106	0.532
78-11-5	PETN				
99-99-0	p-Nitrotoluene	.16	U	0.160	0.532
99-99-0	p-Nitrotoluene				
3058-38-6	TATB	.319	U	0.319	1.06
3058-38-6	TATB				
618-87-1	3,5-Dinitroaniline	.319	U	0.319	1.06
618-87-1	3,5-Dinitroaniline				
78-30-8	tris(o-cresyl) phosphate	.319	U	0.319	1.06
78-30-8	tris(o-cresyl) phosphate				
59229-75-3	2,6-Diamino-4-nitrotoluene	.532	U	0.532	2.66
59229-75-3	2,6-Diamino-4-nitrotoluene				
6629-29-4	2,4-Diamino-6-nitrotoluene	.532	U	0.532	2.66
6629-29-4	2,4-Diamino-6-nitrotoluene				

Quality Control Summary

High Explosives Surrogate Recovery Summary

Lab Name: GEL Laboratories LLC

GEL Job No (SDG): 2018-1811

Lab Code: GEL

HPLC Column: Ultracarb Phenomenex 5u ODS (20)

Lab Sample ID	Client Sample ID	DNT	QC Limits	Flg
444566003	CAPA-18-2	95	55 - 115	
444566007	CAPA-18-12	106	55 - 115	
444566008	CAWA-18-54	96	55 - 115	
444566013	CAWA-18-56	99	55 - 115	
1203979520	MB for batch 1742520	93	55 - 115	
1203979521	LCS for batch 1742520	80	55 - 115	
1203979522	CAPA-18-2MS	89	55 - 115	
1203979523	CAPA-18-2MSD	87	55 - 115	

DNT = 3,4-Dinitrotoluene

3B
High Explosives LCS/LCS Duplicate Summary

Lab Name: GEL Laboratories LLC

Client ID: LCS

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Extract Batch Code: 1742520

Date Extracted: 27-FEB-18

GEL LCS ID: 1203979521

GEL LCSDUP ID: .

Analysis Date/Time: 05-MAR-18 19:33

DUP Analysis Date/Time:

Reporting Units: ug/L

QC Type: LCS/LCSD

Compound	Spike Added	LCS Conc	LCS Rec #	LCSD Conc	LCSD Rec #	RPD #	RPD	Recovery Limits
Tetryl	5	3.54	71					55 - 122
m-Dinitrobenzene	5	4.77	95					74 - 117
m-Nitrotoluene	5	3.94	79					66 - 114
o-Nitrotoluene	5	4.22	84					64 - 115
p-Nitrotoluene	5	4.35	87					66 - 127
tris(o-cresyl) phosphate	5	3.6	72					43 - 104
1,3,5-Trinitrobenzene	5	4.93	99					70 - 110
2,4,6-Trinitrotoluene	5	4.19	84					69 - 113
2,4-Diamino-6-nitrotoluene	5	5.16	103					50 - 121
2,4-Dinitrotoluene	5	4.01	80					71 - 110
2,6-Diamino-4-nitrotoluene	5	4.55	91					53 - 127
2,6-Dinitrotoluene	5	4.39	88					72 - 105
2-Amino-4,6-dinitrotoluene	5	4.29	86					70 - 112
3,5-Dinitroaniline	5	4.19	84					70 - 121
4-Amino-2,6-dinitrotoluene	5	4.33	87					74 - 116
DNX	5	4.9	98					65 - 113
HMX	5	4.71	94					58 - 113
MXN	5	5.4	108					66 - 114
Nitrobenzene	5	4.95	99					64 - 115
PETN	5	5.04	101					57 - 126
RDX	5	4.48	90					64 - 117
TATB	5	4.1	82					47 - 135
TNX	5	4.56	91					51 - 110

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

3
High Explosives MS/MSD Summary

Lab Name: GEL Laboratories LLC

Client ID: CAPA-18-2

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Extract Batch Code: 1742520

Date Extracted: 27-FEB-18

GEL Spike ID: 1203979522

GEL SpikeDup ID: 1203979523

Analysis Date/Time: 05-MAR-18 20:44

MSD Analysis Date/Time: 05-MAR-18 21:20

Reporting Units: ug/L

QC Type: MS/MSD

Compound	Spike Added	Sample Conc	MS Conc	MS Rec #	MSD Conc	MSD Rec #	RPD #	RPD Limit	Rec Limits
p-Nitrotoluene	5.81395	0	5.39	93	5.01	87	7	30	61 - 129
tris(o-cresyl) phosphate	5.81395	0	4.24	73	3.8	66	11	30	38 - 105
1,3,5-Trinitrobenzene	5.81395	0	5.44	94	5.72	100	5	30	67 - 111
2,4,6-Trinitrotoluene	5.81395	0	5.22	90	4.74	82	10	30	66 - 112
2,4-Diamino-6-nitrotoluene	5.81395	0	6.51	112	6.02	105	8	30	50 - 121
2,4-Dinitrotoluene	5.81395	0	5.82	100	5.38	94	8	30	69 - 113
2,6-Diamino-4-nitrotoluene	5.81395	0	5.51	95	5.61	98	2	30	53 - 127
2,6-Dinitrotoluene	5.81395	0	5.18	89	4.85	84	7	30	70 - 106
2-Amino-4,6-dinitrotoluene	5.81395	.12	5.79	98	5.2	88	11	30	67 - 115
3,5-Dinitroaniline	5.81395	.0492	5.45	93	5.03	87	8	30	70 - 121
4-Amino-2,6-dinitrotoluene	5.81395	.195	5.81	97	5.45	91	6	30	65 - 120
DNX	5.81395	0	5.23	90	5.72	100	9	30	53 - 124
HMX	5.81395	6.64	12.8	106	12.8	107	0	30	44 - 128
MNX	5.81395	.0658	5.3	90	5.97	103	12	30	60 - 121
Nitrobenzene	5.81395	0	5.71	98	5.86	102	2	30	62 - 116
PETN	5.81395	0	6.52	112	5.93	103	9	30	51 - 131
RDX	5.81395	7.69	12.7	86	13.1	94	3	30	57 - 125
TATB	5.81395	.0341	4.94	84	5.13	89	4	30	38 - 149
TNX	5.81395	0	5.28	91	5.46	95	3	30	46 - 120
Tetryl	5.81395	0	4.17	72	4.2	73	1	30	50 - 126
m-Dinitrobenzene	5.81395	0	5.88	101	5.54	96	6	30	74 - 117
m-Nitrotoluene	5.81395	0	5.31	91	5.07	88	5	30	59 - 120
o-Nitrotoluene	5.81395	0	5.19	89	4.66	81	11	30	56 - 119

#Column to be used to flag recovery and RPD values with an asterisk

Quality Control Data

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: MB for batch 1742520

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 1203979520

Sample Amount 1000 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

GEL data file: EXP0305016.wiff

Date Analyzed: 05-MAR-18 18:58

Dilution Factor: 2

Concentration Units: ug/L

Cas No.	Compound	Concentration*	Q	MDL	PQL
118-96-7	2,4,6-Trinitrotoluene	.08	U	0.080	0.250
<i>118-96-7</i>	<i>2,4,6-Trinitrotoluene</i>				
121-14-2	2,4-Dinitrotoluene	.08	U	0.080	0.250
<i>121-14-2</i>	<i>2,4-Dinitrotoluene</i>				
121-82-4	RDX	.08	U	0.080	0.250
<i>121-82-4</i>	<i>RDX</i>				
13980-04-6	TNX	.08	U	0.080	0.250
<i>13980-04-6</i>	<i>TNX</i>				
19406-51-0	4-Amino-2,6-dinitrotoluene	.08	U	0.080	0.250
<i>19406-51-0</i>	<i>4-Amino-2,6-dinitrotoluene</i>				
2691-41-0	HMX	.08	U	0.080	0.250
<i>2691-41-0</i>	<i>HMX</i>				
35572-78-2	2-Amino-4,6-dinitrotoluene	.08	U	0.080	0.250
<i>35572-78-2</i>	<i>2-Amino-4,6-dinitrotoluene</i>				
479-45-8	Tetryl	.08	U	0.080	0.500
<i>479-45-8</i>	<i>Tetryl</i>				
5755-27-1	MNX	.08	U	0.080	0.250
<i>5755-27-1</i>	<i>MNX</i>				
606-20-2	2,6-Dinitrotoluene	.08	U	0.080	0.250
<i>606-20-2</i>	<i>2,6-Dinitrotoluene</i>				
80251-29-2	DNX	.08	U	0.080	0.250
<i>80251-29-2</i>	<i>DNX</i>				
98-95-3	Nitrobenzene	.08	U	0.080	0.250
<i>98-95-3</i>	<i>Nitrobenzene</i>				
99-08-1	m-Nitrotoluene	.08	U	0.080	0.250
<i>99-08-1</i>	<i>m-Nitrotoluene</i>				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: MB for batch 1742520

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 1203979520

Sample Amount 1000 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

Cas No.	Compound	Concentration*	Q	MDL	PQL
99-35-4	1,3,5-Trinitrobenzene	.08	U	0.080	0.250
99-35-4	1,3,5-Trinitrobenzene				
99-65-0	m-Dinitrobenzene	.08	U	0.080	0.250
99-65-0	m-Dinitrobenzene				
88-72-2	o-Nitrotoluene	.082	U	0.082	0.250
88-72-2	o-Nitrotoluene				
78-11-5	PETN	.1	U	0.100	0.500
78-11-5	PETN				
99-99-0	p-Nitrotoluene	.15	U	0.150	0.500
99-99-0	p-Nitrotoluene				
3058-38-6	TATB	.3	U	0.300	1.00
3058-38-6	TATB				
618-87-1	3,5-Dinitroaniline	.3	U	0.300	1.00
618-87-1	3,5-Dinitroaniline				
78-30-8	tris(o-cresyl) phosphate	.3	U	0.300	1.00
78-30-8	tris(o-cresyl) phosphate				
59229-75-3	2,6-Diamino-4-nitrotoluene	.5	U	0.500	2.50
59229-75-3	2,6-Diamino-4-nitrotoluene				
6629-29-4	2,4-Diamino-6-nitrotoluene	.5	U	0.500	2.50
6629-29-4	2,4-Diamino-6-nitrotoluene				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: LCS for batch 1742520

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 1203979521

Sample Amount 1000 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

GEL data file: EXP0305017.wiff

Date Analyzed: 05-MAR-18 19:33

Dilution Factor: 2

Concentration Units: ug/L

Cas No.	Compound	Concentration*	Q	MDL	PQL
479-45-8	Tetryl	3.54		0.080	0.500
<i>479-45-8</i>	<i>Tetryl</i>				
78-30-8	tris(o-cresyl) phosphate	3.6		0.300	1.00
<i>78-30-8</i>	<i>tris(o-cresyl) phosphate</i>				
99-08-1	m-Nitrotoluene	3.94		0.080	0.250
<i>99-08-1</i>	<i>m-Nitrotoluene</i>				
121-14-2	2,4-Dinitrotoluene	4.01		0.080	0.250
<i>121-14-2</i>	<i>2,4-Dinitrotoluene</i>				
3058-38-6	TATB	4.1		0.300	1.00
<i>3058-38-6</i>	<i>TATB</i>				
118-96-7	2,4,6-Trinitrotoluene	4.19		0.080	0.250
<i>118-96-7</i>	<i>2,4,6-Trinitrotoluene</i>				
618-87-1	3,5-Dinitroaniline	4.19		0.300	1.00
<i>618-87-1</i>	<i>3,5-Dinitroaniline</i>				
88-72-2	o-Nitrotoluene	4.22		0.082	0.250
<i>88-72-2</i>	<i>o-Nitrotoluene</i>				
35572-78-2	2-Amino-4,6-dinitrotoluene	4.29		0.080	0.250
<i>35572-78-2</i>	<i>2-Amino-4,6-dinitrotoluene</i>				
19406-51-0	4-Amino-2,6-dinitrotoluene	4.33		0.080	0.250
<i>19406-51-0</i>	<i>4-Amino-2,6-dinitrotoluene</i>				
99-99-0	p-Nitrotoluene	4.35		0.150	0.500
<i>99-99-0</i>	<i>p-Nitrotoluene</i>				
606-20-2	2,6-Dinitrotoluene	4.39		0.080	0.250
<i>606-20-2</i>	<i>2,6-Dinitrotoluene</i>				
121-82-4	RDX	4.48		0.080	0.250
<i>121-82-4</i>	<i>RDX</i>				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: LCS for batch 1742520

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 1203979521

Sample Amount 1000 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

Cas No.	Compound	Concentration*	Q	MDL	PQL
59229-75-3	2,6-Diamino-4-nitrotoluene	4.55		0.500	2.50
<i>59229-75-3</i>	<i>2,6-Diamino-4-nitrotoluene</i>				
13980-04-6	TNX	4.56		0.080	0.250
<i>13980-04-6</i>	<i>TNX</i>				
2691-41-0	HMX	4.71		0.080	0.250
<i>2691-41-0</i>	<i>HMX</i>				
99-65-0	m-Dinitrobenzene	4.77		0.080	0.250
<i>99-65-0</i>	<i>m-Dinitrobenzene</i>				
80251-29-2	DNX	4.9		0.080	0.250
<i>80251-29-2</i>	<i>DNX</i>				
99-35-4	1,3,5-Trinitrobenzene	4.93		0.080	0.250
<i>99-35-4</i>	<i>1,3,5-Trinitrobenzene</i>				
98-95-3	Nitrobenzene	4.95		0.080	0.250
<i>98-95-3</i>	<i>Nitrobenzene</i>				
78-11-5	PETN	5.04		0.100	0.500
<i>78-11-5</i>	<i>PETN</i>				
6629-29-4	2,4-Diamino-6-nitrotoluene	5.16		0.500	2.50
<i>6629-29-4</i>	<i>2,4-Diamino-6-nitrotoluene</i>				
5755-27-1	MNX	5.4		0.080	0.250
<i>5755-27-1</i>	<i>MNX</i>				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAPA-18-2(444566003MS)MS

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 1203979522

Sample Amount 860 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

GEL data file: EXP0305019.wiff

Date Analyzed: 05-MAR-18 20:44

Dilution Factor: 2

Concentration Units: ug/L

Cas No.	Compound	Concentration*	Q	MDL	PQL
479-45-8	Tetryl	4.17		0.093	0.581
<i>479-45-8</i>	<i>Tetryl</i>				
78-30-8	tris(o-cresyl) phosphate	4.24		0.349	1.16
<i>78-30-8</i>	<i>tris(o-cresyl) phosphate</i>				
3058-38-6	TATB	4.94		0.349	1.16
<i>3058-38-6</i>	<i>TATB</i>				
606-20-2	2,6-Dinitrotoluene	5.18		0.093	0.291
<i>606-20-2</i>	<i>2,6-Dinitrotoluene</i>				
88-72-2	o-Nitrotoluene	5.19		0.0953	0.291
<i>88-72-2</i>	<i>o-Nitrotoluene</i>				
118-96-7	2,4,6-Trinitrotoluene	5.22		0.093	0.291
<i>118-96-7</i>	<i>2,4,6-Trinitrotoluene</i>				
80251-29-2	DNX	5.23		0.093	0.291
<i>80251-29-2</i>	<i>DNX</i>				
13980-04-6	TNX	5.28		0.093	0.291
<i>13980-04-6</i>	<i>TNX</i>				
5755-27-1	MNX	5.3		0.093	0.291
<i>5755-27-1</i>	<i>MNX</i>				
99-08-1	m-Nitrotoluene	5.31		0.093	0.291
<i>99-08-1</i>	<i>m-Nitrotoluene</i>				
99-99-0	p-Nitrotoluene	5.39		0.174	0.581
<i>99-99-0</i>	<i>p-Nitrotoluene</i>				
99-35-4	1,3,5-Trinitrobenzene	5.44		0.093	0.291
<i>99-35-4</i>	<i>1,3,5-Trinitrobenzene</i>				
618-87-1	3,5-Dinitroaniline	5.45		0.349	1.16
<i>618-87-1</i>	<i>3,5-Dinitroaniline</i>				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAPA-18-2(444566003MS)MS

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 1203979522

Sample Amount 860 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

Cas No.	Compound	Concentration*	Q	MDL	PQL
59229-75-3	2,6-Diamino-4-nitrotoluene	5.51		0.581	2.91
<i>59229-75-3</i>	<i>2,6-Diamino-4-nitrotoluene</i>				
98-95-3	Nitrobenzene	5.71		0.093	0.291
<i>98-95-3</i>	<i>Nitrobenzene</i>				
35572-78-2	2-Amino-4,6-dinitrotoluene	5.79		0.093	0.291
<i>35572-78-2</i>	<i>2-Amino-4,6-dinitrotoluene</i>				
19406-51-0	4-Amino-2,6-dinitrotoluene	5.81		0.093	0.291
<i>19406-51-0</i>	<i>4-Amino-2,6-dinitrotoluene</i>				
121-14-2	2,4-Dinitrotoluene	5.82		0.093	0.291
<i>121-14-2</i>	<i>2,4-Dinitrotoluene</i>				
99-65-0	m-Dinitrobenzene	5.88		0.093	0.291
<i>99-65-0</i>	<i>m-Dinitrobenzene</i>				
6629-29-4	2,4-Diamino-6-nitrotoluene	6.51		0.581	2.91
<i>6629-29-4</i>	<i>2,4-Diamino-6-nitrotoluene</i>				
78-11-5	PETN	6.52		0.116	0.581
<i>78-11-5</i>	<i>PETN</i>				
121-82-4	RDX	12.7		0.093	0.291
<i>121-82-4</i>	<i>RDX</i>				
2691-41-0	HMX	12.8		0.093	0.291
<i>2691-41-0</i>	<i>HMX</i>				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAPA-18-2(444566003MSD)MSD

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 1203979523

Sample Amount 870 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

GEL data file: EXP0305020.wiff

Date Analyzed: 05-MAR-18 21:20

Dilution Factor: 2

Concentration Units: ug/L

Cas No.	Compound	Concentration*	Q	MDL	PQL
78-30-8	tris(o-cresyl) phosphate	3.8		0.345	1.15
78-30-8	tris(o-cresyl) phosphate				
479-45-8	Tetryl	4.2		0.092	0.575
479-45-8	Tetryl				
88-72-2	o-Nitrotoluene	4.66		0.0943	0.287
88-72-2	o-Nitrotoluene				
118-96-7	2,4,6-Trinitrotoluene	4.74		0.092	0.287
118-96-7	2,4,6-Trinitrotoluene				
606-20-2	2,6-Dinitrotoluene	4.85		0.092	0.287
606-20-2	2,6-Dinitrotoluene				
99-99-0	p-Nitrotoluene	5.01		0.172	0.575
99-99-0	p-Nitrotoluene				
618-87-1	3,5-Dinitroaniline	5.03		0.345	1.15
618-87-1	3,5-Dinitroaniline				
99-08-1	m-Nitrotoluene	5.07		0.092	0.287
99-08-1	m-Nitrotoluene				
3058-38-6	TATB	5.13		0.345	1.15
3058-38-6	TATB				
35572-78-2	2-Amino-4,6-dinitrotoluene	5.2		0.092	0.287
35572-78-2	2-Amino-4,6-dinitrotoluene				
121-14-2	2,4-Dinitrotoluene	5.38		0.092	0.287
121-14-2	2,4-Dinitrotoluene				
19406-51-0	4-Amino-2,6-dinitrotoluene	5.45		0.092	0.287
19406-51-0	4-Amino-2,6-dinitrotoluene				
13980-04-6	TNX	5.46		0.092	0.287
13980-04-6	TNX				

1
High Explosives Analysis Data Sheet

Lab Name: GEL Laboratories LLC

Client Sample ID: CAPA-18-2(444566003MSD)MSD

Lab Code: GEL

GEL Job No (SDG) 2018-1811

Matrix: WATER

GEL Sample ID: 1203979523

Sample Amount 870 mL

Date Received: 23-FEB-18

Moisture: .

Extraction Batch ID: 1742520

Extraction Type Sol Exchange

Date Extracted: 27-FEB-18

Concentrated Extract Volume (mL) 5

Injection Volume (uL):50

Cas No.	Compound	Concentration*	Q	MDL	PQL
99-65-0	m-Dinitrobenzene	5.54		0.092	0.287
99-65-0	<i>m-Dinitrobenzene</i>				
59229-75-3	2,6-Diamino-4-nitrotoluene	5.61		0.575	2.87
59229-75-3	<i>2,6-Diamino-4-nitrotoluene</i>				
80251-29-2	DNX	5.72		0.092	0.287
80251-29-2	<i>DNX</i>				
99-35-4	1,3,5-Trinitrobenzene	5.72		0.092	0.287
99-35-4	<i>1,3,5-Trinitrobenzene</i>				
98-95-3	Nitrobenzene	5.86		0.092	0.287
98-95-3	<i>Nitrobenzene</i>				
78-11-5	PETN	5.93		0.115	0.575
78-11-5	<i>PETN</i>				
5755-27-1	MNX	5.97		0.092	0.287
5755-27-1	<i>MNX</i>				
6629-29-4	2,4-Diamino-6-nitrotoluene	6.02		0.575	2.87
6629-29-4	<i>2,4-Diamino-6-nitrotoluene</i>				
2691-41-0	HMX	12.8		0.092	0.287
2691-41-0	<i>HMX</i>				
121-82-4	RDX	13.1		0.092	0.287
121-82-4	<i>RDX</i>				

Explosives Initial Calibration Blank

Lab Name: GEL Laboratories LLCGEL Job No(SDG): 2018-1811Lab Code: GELLab Sample ID: XIBLK01Analysis Date: 05-MAR-18 10:05GEL Data File: EXP0305001.wiffInstrument ID: LCMSMS5Column: Ultracarb Phenomenex 5u ODS (20)

Compound	True	Found (ug/L)
3,4-Dinitrotoluene	0	0
tris(o-cresyl) phosphate	0	0
TATB	0	0
3,5-Dinitroaniline	0	0
2,4-Diamino-6-nitrotoluene	0	0
2,6-Diamino-4-nitrotoluene	0	0
DNX	0	0
MNX	0	0
TNX	0	0
1,3,5-Trinitrobenzene	0	0
2,4,6-Trinitrotoluene	0	0
2,4-Dinitrotoluene	0	0
2,6-Dinitrotoluene	0	0
2-Amino-4,6-dinitrotoluene	0	0
4-Amino-2,6-dinitrotoluene	0	0
HMX	0	0
Nitrobenzene	0	0
PETN	0	0
RDX	0	0
Tetryl	0	0
m-Dinitrobenzene	0	0
m-Nitrotoluene	0	0
o-Nitrotoluene	0	0
p-Nitrotoluene	0	0

Explosives Initial Calibration Blank

Lab Name: GEL Laboratories LLCGEL Job No(SDG): 2018-1811Lab Code: GELLab Sample ID: XIBLK01Analysis Date: 05-MAR-18 10:40GEL Data File: EXP0305002.wiffInstrument ID: LCMSMS5Column: Ultracarb Phenomenex 5u ODS (20)

Compound	True	Found (ug/L)
TNX	0	0
1,3,5-Trinitrobenzene	0	0
2,4,6-Trinitrotoluene	0	0
2,4-Dinitrotoluene	0	0
2,6-Dinitrotoluene	0	0
2-Amino-4,6-dinitrotoluene	0	0
4-Amino-2,6-dinitrotoluene	0	0
HMX	0	0
Nitrobenzene	0	0
PETN	0	0
RDX	0	0
Tetryl	0	0
m-Dinitrobenzene	0	0
m-Nitrotoluene	0	0
o-Nitrotoluene	0	0
p-Nitrotoluene	0	0
MNX	0	0
3,4-Dinitrotoluene	0	0
tris(o-cresyl) phosphate	0	0
TATB	0	0
3,5-Dinitroaniline	0	0
2,4-Diamino-6-nitrotoluene	0	0
2,6-Diamino-4-nitrotoluene	0	0
DNX	0	0

4A
Explosives Continuing Calibration Blank

Lab Name: GEL Laboratories LLC

GEL Job No(SDG): 2018-1811

Lab Code: GEL

Lab Sample ID: XIBLK02

Analysis Date: 05-MAR-18 15:25

GEL Data File: EXP0305010.wiff

Instrument ID: LCMSMS5

Column: Ultracarb Phenomenex 5u ODS (20)

Compound	True	Found (ug/L)
3,4-Dinitrotoluene	0	0
tris(o-cresyl) phosphate	0	0
TATB	0	0
3,5-Dinitroaniline	0	0
2,4-Diamino-6-nitrotoluene	0	0
2,6-Diamino-4-nitrotoluene	0	0
DNX	0	0
MNX	0	0
TNX	0	0
1,3,5-Trinitrobenzene	0	0
2,4,6-Trinitrotoluene	0	0
2,4-Dinitrotoluene	0	0
2,6-Dinitrotoluene	0	0
2-Amino-4,6-dinitrotoluene	0	0
4-Amino-2,6-dinitrotoluene	0	0
HMX	0	0
Nitrobenzene	0	0
PETN	0	0
RDX	0	0
Tetryl	0	0
m-Dinitrobenzene	0	0
m-Nitrotoluene	0	0
o-Nitrotoluene	0	0
p-Nitrotoluene	0	0

4A
Explosives Continuing Calibration Blank

Lab Name: GEL Laboratories LLC

GEL Job No(SDG): 2018-1811

Lab Code: GEL

Lab Sample ID: XIBLK03

Analysis Date: 05-MAR-18 17:47

GEL Data File: EXP0305014.wiff

Instrument ID: LCMSMS5

Column: Ultracarb Phenomenex 5u ODS (20)

Compound	True	Found (ug/L)
2,4-Diamino-6-nitrotoluene	0	0
2,6-Diamino-4-nitrotoluene	0	0
DNX	0	0
MNX	0	0
TNX	0	0
1,3,5-Trinitrobenzene	0	0
2,4,6-Trinitrotoluene	0	0
2,4-Dinitrotoluene	0	0
2,6-Dinitrotoluene	0	0
2-Amino-4,6-dinitrotoluene	0	0
4-Amino-2,6-dinitrotoluene	0	0
HMX	0	0
Nitrobenzene	0	0
PETN	0	0
RDX	0	0
Tetryl	0	0
m-Dinitrobenzene	0	0
m-Nitrotoluene	0	0
o-Nitrotoluene	0	0
p-Nitrotoluene	0	0
3,4-Dinitrotoluene	0	0
tris(o-cresyl) phosphate	0	0
TATB	0	0
3,5-Dinitroaniline	0	0

4A
Explosives Continuing Calibration Blank

Lab Name: GEL Laboratories LLC

GEL Job No(SDG): 2018-1811

Lab Code: GEL

Lab Sample ID: XIBLK04

Analysis Date: 05-MAR-18 21:56

GEL Data File: EXP0305021.wiff

Instrument ID: LCMSMS5

Column: Ultracarb Phenomenex 5u ODS (20)

Compound	True	Found (ug/L)
3,4-Dinitrotoluene	0	0
tris(o-cresyl) phosphate	0	0
TATB	0	0
3,5-Dinitroaniline	0	0
2,4-Diamino-6-nitrotoluene	0	0
2,6-Diamino-4-nitrotoluene	0	0
DNX	0	0
MNX	0	0
TNX	0	0
1,3,5-Trinitrobenzene	0	0
2,4,6-Trinitrotoluene	0	0
2,4-Dinitrotoluene	0	0
2,6-Dinitrotoluene	0	0
2-Amino-4,6-dinitrotoluene	0	0
4-Amino-2,6-dinitrotoluene	0	0
HMX	0	0
Nitrobenzene	0	0
PETN	0	0
RDX	0	0
Tetryl	0	0
m-Dinitrobenzene	0	0
m-Nitrotoluene	0	0
o-Nitrotoluene	0	0
p-Nitrotoluene	0	0

4A
Explosives Continuing Calibration Blank

Lab Name: GEL Laboratories LLC

GEL Job No(SDG): 2018-1811

Lab Code: GEL

Lab Sample ID: XIBLK05

Analysis Date: 06-MAR-18 00:18

GEL Data File: EXP0305025.wiff

Instrument ID: LCMSMS5

Column: Ultracarb Phenomenex 5u ODS (20)

Compound	True	Found (ug/L)
tris(o-cresyl) phosphate	0	0
TATB	0	0
3,5-Dinitroaniline	0	0
2,4-Diamino-6-nitrotoluene	0	0
2,6-Diamino-4-nitrotoluene	0	0
DNX	0	0
MNX	0	0
TNX	0	0
1,3,5-Trinitrobenzene	0	0
2,4,6-Trinitrotoluene	0	0
2,4-Dinitrotoluene	0	0
2,6-Dinitrotoluene	0	0
2-Amino-4,6-dinitrotoluene	0	0
4-Amino-2,6-dinitrotoluene	0	0
HMX	0	0
Nitrobenzene	0	0
PETN	0	0
RDX	0	0
Tetryl	0	0
m-Dinitrobenzene	0	0
m-Nitrotoluene	0	0
o-Nitrotoluene	0	0
p-Nitrotoluene	0	0
3,4-Dinitrotoluene	0	0

4A
Explosives Continuing Calibration Blank

Lab Name: GEL Laboratories LLC

GEL Job No(SDG): 2018-1811

Lab Code: GEL

Lab Sample ID: XIBLK06

Analysis Date: 06-MAR-18 01:29

GEL Data File: EXP0305027.wiff

Instrument ID: LCMSMS5

Column: Ultracarb Phenomenex 5u ODS (20)

Compound	True	Found (ug/L)
o-Nitrotoluene	0	0
p-Nitrotoluene	0	0
3,4-Dinitrotoluene	0	0
tris(o-cresyl) phosphate	0	0
TATB	0	0
3,5-Dinitroaniline	0	0
2,4-Diamino-6-nitrotoluene	0	0
2,6-Diamino-4-nitrotoluene	0	0
DNX	0	0
MXN	0	0
TNX	0	0
1,3,5-Trinitrobenzene	0	0
2,4,6-Trinitrotoluene	0	0
2,4-Dinitrotoluene	0	0
2,6-Dinitrotoluene	0	0
2-Amino-4,6-dinitrotoluene	0	0
4-Amino-2,6-dinitrotoluene	0	0
HMX	0	0
Nitrobenzene	0	0
PETN	0	0
RDX	0	0
Tetryl	0	0
m-Dinitrobenzene	0	0
m-Nitrotoluene	0	0

Metals Analysis

Case Narrative

Metals
Technical Case Narrative
ARS International, LLC (ARSL)
SDG #: 2018-1811
Work Order #: 444566

Sample ID	Client ID
444566001	CAPA-18-1
444566002	CAPA-18-2
444566005	CAPA-18-10
444566006	CAPA-18-12
444566011	CAWA-18-55
444566012	CAWA-18-56
1203978044	Method Blank (MB) ICP
1203978045	Laboratory Control Sample (LCS)
1203978048	444566001(CAPA-18-1L) Serial Dilution (SD)
1203978046	444566001(CAPA-18-1D) Sample Duplicate (DUP)
1203978047	444566001(CAPA-18-1S) Matrix Spike (MS)
1203978054	Method Blank (MB) ICP-MS
1203978055	Laboratory Control Sample (LCS)
1203978058	444566001(CAPA-18-1L) Serial Dilution (SD)
1203978056	444566001(CAPA-18-1D) Sample Duplicate (DUP)
1203978057	444566001(CAPA-18-1S) Matrix Spike (MS)
1203979604	Method Blank (MB) CVAA
1203979605	Laboratory Control Sample (LCS)
1203979608	444566001(CAPA-18-1L) Serial Dilution (SD)
1203979606	444566001(CAPA-18-1D) Sample Duplicate (DUP)
1203979607	444566001(CAPA-18-1S) Matrix Spike (MS)

Sample Analysis

Samples 444566001,002,005,006,011 and 012 in this SDG were analyzed for metals and mercury on an "as received" basis.

Method/Analysis Information

Analytical Batch:	1741857, 1741864, 1742549 and 1747180
Prep Batch :	1741856, 1741863 and 1742548
Standard Operating Procedures:	GL-MA-E-013 REV# 30, GL-MA-E-006 REV# 14, GL-MA-E-014 REV# 32, GL-MA-E-010 REV# 36 and GL-GC-E-107 REV# 10
Analytical Method:	SW846 3005A/6010C, SW846 3005A/6020A, EPA 245.2 1974 and SM:A2340B
Prep Method :	SW846 3005A and EPA 245.1/245.2 Prep

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

System Configuration

The Hardness as CaCO₃ is calculated from Calcium and Magnesium results.

The Metals analysis-ICP was performed on a PE 7300 Optima radial/axial-viewing inductively coupled plasma atomic emission spectrometer. The instrument is equipped with an ESI SC-FAST introduction, cyclonic spray chamber, and yttrium or scandium internal standard.

The Metals analysis-Mercury was performed on a Perkin-Elmer Flow Injection Mercury System (FIMS-100) automated mercury analyzer. The instrument consists of a cold vapor atomic absorption spectrometer set to detect mercury at a wavelength of 253.7 nm.

The Metals analysis - ICPMS was performed on a PerkinElmer NexION 350X ICPMS. The instrument is equipped with a ESI PFA-ST nebulizer, quadrupole mass spectrometer, dual mode electron multiplier detector, and Kinetic Energy Discrimination (KED) technology. Internal standards of scandium, germanium, indium, tantalum, and/or lutetium were utilized to cover the mass spectrum.

Calibration Information

Instrument Calibration

All initial calibration requirements have been met for this sample delivery group (SDG).

CRDL/PQL Requirements

The CRDL/PQL standard recoveries met the referenced advisory control limits.

ICSA/ICSAB Statement

All interference check samples (ICSA and ICSAB) associated with this SDG met the established acceptance criteria. For the ICP-MS analysis, the ICSA solution contains analyte concentrations which are verified trace impurities indigenous to the purchased standard.

Continuing Calibration Blanks (CCB) Requirements

All continuing calibration blanks (CCB) bracketing this batch met the established acceptance criteria.

Continuing Calibration Verification (CCV) Requirements

All continuing calibration verifications (CCV) bracketing this SDG met the acceptance criteria.

Quality Control (QC) Information

Method Blank (MB) Statement

The MBs analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries met the acceptance limits.

Quality Control (QC) Sample Statement

The following samples were selected as the quality control (QC) samples for this SDG: 444566001 (CAPA-18-1)-ICP, ICP-MS and CVAA.

Matrix Spike (MS/MSD) Recovery Statement

The percent recoveries (%R) obtained from the MS/MSD analyses are evaluated when the sample concentration is less than four times (4X) the spike concentration added. The matrix spike met the recommended quality control acceptance criteria for percent recoveries for all applicable analytes.

Duplicate Relative Percent Difference (RPD) Statement

The RPD obtained from the designated sample duplicate (DUP) is evaluated based on acceptance criteria of 20% when the sample is >5X the contract required reporting limit (RL). In cases where either the sample or duplicate value is less than 5X the RL, a control of +/-RL is used to evaluate the DUP results. Not all the applicable analyte RPD values were within the acceptance criteria.

Sample	Analyte	Value
1203978046 (CAPA-18-1DUP)	Tin	abs(18.2 - -1.1)* (+/-10 ug/L)

Serial Dilution % Difference Statement

All applicable analytes in the serial dilution (SDILT) demonstrated acceptable correlation to its associated sample and met the established acceptance percent difference criteria.

Technical Information

Holding Time Specifications

GEL assigns holding times based on the associated methodology. Holding time is measured by comparison of the date and time of sample collection to the date and time of sample preparation and analysis. Those holding times expressed in hours are calculated in the AlphaLIMS system. Those holding times expressed as days expire at midnight on the day of expiration. All samples in this SDG met the specified holding time.

Preparation/Analytical Method Verification

All procedures were performed as stated in the SOP.

Sample Dilutions

The samples in this SDG did not require dilutions.

Preparation Information

The samples in this SDG were not diluted and were prepared according to the cited SOP.

Miscellaneous Information

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. An electronic signature page inserted after the case narrative will include the data validator's signature and title. The signature page also includes the data qualifiers used in the fractional package. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Additional Comments

Total Hardness by Calculation is determined using the results of Total Calcium (Ca) and Total Magnesium (Mg) determined by ICP or ICP-MS.

Hardness = 2.497 (Ca) + 4.118 (Mg)

Please refer to the Total Ca and Total Mg data to validate results appearing on the Hardness Summary sheet. Both results are in the Inorganic/metals section of the package. There is no Batch QC for calculated results, and thus no QC Summary for the Hardness by Calculation Batch. The MDLs and PQLs are calculated using the higher of the two calculated values of Ca or Mg.

Certification Statement

Where the analytical method has been performed under NELAP certification, the analysis has met all of the requirements of the NELAC standard unless otherwise noted in the analytical case narrative.

GEL LABORATORIES LLC

2040 Savage Road Charleston SC 29407 - (843) 556-8171 - www.gel.com

Qualifier Definition Report for

ARSL004 ARS International, LLC (ARS-LANS-MTOA6-25093-GEL)

Client SDG: 2018-1811 GEL Work Order: 444566

The Qualifiers in this report are defined as follows:

- * A quality control analyte recovery is outside of specified acceptance criteria
- J Value is estimated
- U Analyte was analyzed for, but not detected above the MDL, MDA, MDC or LOD.

Review/Validation

GEL requires all analytical data to be verified by a qualified data reviewer. In addition, all CLP-like deliverables receive a third level review of the fractional data package.

The following data validator verified the information presented in this data report:

Signature:



Name: Nik-Cole Elmore

Date: 20 MAR 2018

Title: Data Validator

Sample Data Summary

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811**CONTRACT:** ESHL00114**METHOD TYPE:** EPA**SAMPLE ID:** 444566001**BASIS:** As Received**DATE COLLECTED** 21-FEB-18**CLIENT ID:** CAPA-18-1**LEVEL:** Low**DATE RECEIVED** 23-FEB-18**MATRIX:** W**%SOLIDS:** 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
7439-97-6	Mercury	0.067	ug/L	U	0.067	0.2	0.2	1	AV	MTM1	02/28/18 10:20	022818W1-3	1742549

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811

CONTRACT: ESHL00114

METHOD TYPE: SW846

SAMPLE ID: 444566001

BASIS: As Received

DATE COLLECTED 21-FEB-18

CLIENT ID: CAPA-18-1

LEVEL: Low

DATE RECEIVED 23-FEB-18

MATRIX: W

%SOLIDS: 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
7429-90-5	Aluminum	109	ug/L	J	68	200	200	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-36-0	Antimony	1	ug/L	U	1	3	3	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7440-38-2	Arsenic	2	ug/L	U	2	5	5	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7440-39-3	Barium	75.4	ug/L		1	5	5	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-41-7	Beryllium	1	ug/L	U	1	5	5	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-42-8	Boron	15	ug/L	U	15	50	50	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-43-9	Cadmium	0.30	ug/L	U	0.3	1	1	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7440-70-2	Calcium	23000	ug/L		50	200	200	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-47-3	Chromium	3	ug/L	U	3	10	10	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7440-48-4	Cobalt	1	ug/L	U	1	5	5	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-50-8	Copper	3	ug/L	U	3	10	10	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7439-89-6	Iron	60	ug/L	J	30	100	100	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7439-92-1	Lead	0.50	ug/L	U	0.5	2	2	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7439-95-4	Magnesium	6040	ug/L		110	300	300	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7439-96-5	Manganese	2	ug/L	U	2	10	10	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7439-98-7	Molybdenum	1.14	ug/L		0.2	0.5	0.5	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7440-02-0	Nickel	0.60	ug/L	U	0.6	2	2	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7440-09-7	Potassium	3330	ug/L		50	150	150	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7782-49-2	Selenium	2	ug/L	U	2	5	5	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7631-86-9	Silica	39800	ug/L		53	213	213	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-22-4	Silver	0.30	ug/L	U	0.3	1	1	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7440-23-5	Sodium	24300	ug/L		100	300	300	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-24-6	Strontium	131	ug/L		1	5	5	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-28-0	Thallium	0.60	ug/L	U	0.6	2	2	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7440-31-5	Tin	2.5	ug/L	U*	2.5	10	10	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-61-1	Uranium	0.494	ug/L		0.067	0.2	0.2	1	MS	BAJ	02/27/18 20:57	180227-2	1741864
7440-62-2	Vanadium	2.54	ug/L	J	1	5	5	1	P	TXT1	03/07/18 19:17	030718-1	1741857
7440-66-6	Zinc	3.3	ug/L	U	3.3	10	10	1	P	TXT1	03/07/18 19:17	030718-1	1741857

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811**CONTRACT:** ESHL00114**METHOD TYPE:****SAMPLE ID:** 444566001**BASIS:** As Received**DATE COLLECTED** 21-FEB-18**CLIENT ID:** CAPA-18-1**LEVEL:** Low**DATE RECEIVED** 23-FEB-18**MATRIX:** W**%SOLIDS:** 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
	Hardness as CaCO3	82.2	mg/L		0.453	1.24	1.24	1		TXT1	03/14/18 13:54		1747180

Prep Information:

Analytical Batch	Prep Batch	Prep Method	Initial wt./vol.	Units	Final wt./vol.	Units	Date	Analyst
1741857	1741856	SW846 3005A	50	mL	50	mL	02/23/18	JXM8
1741864	1741863	SW846 3005A	50	mL	50	mL	02/23/18	JXM8
1742549	1742548	EPA 245.1/245.2 Prep	20	mL	20	mL	02/27/18	AXS5

***Analytical Methods:**

P SW846 3005A/6010C
MS SW846 3005A/6020A
AV EPA 245.2 1974

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811**CONTRACT:** ESHL00114**METHOD TYPE:** EPA**SAMPLE ID:** 444566002**BASIS:** As Received**DATE COLLECTED** 21-FEB-18**CLIENT ID:** CAPA-18-2**LEVEL:** Low**DATE RECEIVED** 23-FEB-18**MATRIX:** W**%SOLIDS:** 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
7439-97-6	Mercury	0.067	ug/L	U	0.067	0.2	0.2	1	AV	MTM1	02/28/18 10:28	022818W1-3	1742549

Prep Information:

Analytical Batch	Prep Batch	Prep Method	Initial wt./vol.	Units	Final wt./vol.	Units	Date	Analyst
1742549	1742548	EPA 245.1/245.2 Prep	20	mL	20	mL	02/27/18	AXS5

***Analytical Methods:**

AV EPA 245.2 1974

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811**CONTRACT:** ESHL00114**METHOD TYPE:** EPA**SAMPLE ID:** 444566005**BASIS:** As Received**DATE COLLECTED** 21-FEB-18**CLIENT ID:** CAPA-18-10**LEVEL:** Low**DATE RECEIVED** 23-FEB-18**MATRIX:** W**%SOLIDS:** 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
7439-97-6	Mercury	0.067	ug/L	U	0.067	0.2	0.2	1	AV	MTM1	02/28/18 10:30	022818W1-3	1742549

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811

CONTRACT: ESHL00114

METHOD TYPE: SW846

SAMPLE ID: 444566005

BASIS: As Received

DATE COLLECTED 21-FEB-18

CLIENT ID: CAPA-18-10

LEVEL: Low

DATE RECEIVED 23-FEB-18

MATRIX: W

%SOLIDS: 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
7429-90-5	Aluminum	111	ug/L	J	68	200	200	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-36-0	Antimony	1	ug/L	U	1	3	3	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7440-38-2	Arsenic	2.18	ug/L	J	2	5	5	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7440-39-3	Barium	75.6	ug/L		1	5	5	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-41-7	Beryllium	1	ug/L	U	1	5	5	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-42-8	Boron	15	ug/L	U	15	50	50	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-43-9	Cadmium	0.30	ug/L	U	0.3	1	1	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7440-70-2	Calcium	23100	ug/L		50	200	200	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-47-3	Chromium	3	ug/L	U	3	10	10	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7440-48-4	Cobalt	1	ug/L	U	1	5	5	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-50-8	Copper	3	ug/L	U	3	10	10	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7439-89-6	Iron	66.8	ug/L	J	30	100	100	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7439-92-1	Lead	0.50	ug/L	U	0.5	2	2	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7439-95-4	Magnesium	6130	ug/L		110	300	300	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7439-96-5	Manganese	2	ug/L	U	2	10	10	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7439-98-7	Molybdenum	1.12	ug/L		0.2	0.5	0.5	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7440-02-0	Nickel	0.623	ug/L	J	0.6	2	2	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7440-09-7	Potassium	3330	ug/L		50	150	150	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7782-49-2	Selenium	2	ug/L	U	2	5	5	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7631-86-9	Silica	39500	ug/L		53	213	213	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-22-4	Silver	0.30	ug/L	U	0.3	1	1	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7440-23-5	Sodium	24300	ug/L		100	300	300	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-24-6	Strontium	131	ug/L		1	5	5	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-28-0	Thallium	0.60	ug/L	U	0.6	2	2	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7440-31-5	Tin	2.5	ug/L	U*	2.5	10	10	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-61-1	Uranium	0.502	ug/L		0.067	0.2	0.2	1	MS	BAJ	02/27/18 21:13	180227-2	1741864
7440-62-2	Vanadium	1.8	ug/L	J	1	5	5	1	P	TXT1	03/07/18 19:11	030718-1	1741857
7440-66-6	Zinc	3.3	ug/L	U	3.3	10	10	1	P	TXT1	03/07/18 19:11	030718-1	1741857

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811**CONTRACT:** ESHL00114**METHOD TYPE:****SAMPLE ID:** 444566005**BASIS:** As Received**DATE COLLECTED** 21-FEB-18**CLIENT ID:** CAPA-18-10**LEVEL:** Low**DATE RECEIVED** 23-FEB-18**MATRIX:** W**%SOLIDS:** 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
	Hardness as CaCO3	82.8	mg/L		0.453	1.24	1.24	1		TXT1	03/14/18 13:54		1747180

Prep Information:

Analytical Batch	Prep Batch	Prep Method	Initial wt./vol.	Units	Final wt./vol.	Units	Date	Analyst
1741857	1741856	SW846 3005A	50	mL	50	mL	02/23/18	JXM8
1741864	1741863	SW846 3005A	50	mL	50	mL	02/23/18	JXM8
1742549	1742548	EPA 245.1/245.2 Prep	20	mL	20	mL	02/27/18	AXS5

***Analytical Methods:**

P SW846 3005A/6010C
MS SW846 3005A/6020A
AV EPA 245.2 1974

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811**CONTRACT:** ESHL00114**METHOD TYPE:** EPA**SAMPLE ID:** 444566006**BASIS:** As Received**DATE COLLECTED** 21-FEB-18**CLIENT ID:** CAPA-18-12**LEVEL:** Low**DATE RECEIVED** 23-FEB-18**MATRIX:** W**%SOLIDS:** 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
7439-97-6	Mercury	0.067	ug/L	U	0.067	0.2	0.2	1	AV	MTM1	02/28/18 10:31	022818W1-3	1742549

Prep Information:

Analytical Batch	Prep Batch	Prep Method	Initial wt./vol.	Units	Final wt./vol.	Units	Date	Analyst
1742549	1742548	EPA 245.1/245.2 Prep	20	mL	20	mL	02/27/18	AXS5

***Analytical Methods:**

AV EPA 245.2 1974

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811**CONTRACT:** ESHL00114**METHOD TYPE:** EPA**SAMPLE ID:** 444566011**BASIS:** As Received**DATE COLLECTED** 21-FEB-18**CLIENT ID:** CAWA-18-55**LEVEL:** Low**DATE RECEIVED** 23-FEB-18**MATRIX:** W**%SOLIDS:** 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
7439-97-6	Mercury	0.067	ug/L	U	0.067	0.2	0.2	1	AV	MTM1	02/28/18 10:36	022818W1-3	1742549

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811

CONTRACT: ESHL00114

METHOD TYPE: SW846

SAMPLE ID: 444566011

BASIS: As Received

DATE COLLECTED 21-FEB-18

CLIENT ID: CAWA-18-55

LEVEL: Low

DATE RECEIVED 23-FEB-18

MATRIX: W

%SOLIDS: 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
7429-90-5	Aluminum	68	ug/L	U	68	200	200	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-36-0	Antimony	1	ug/L	U	1	3	3	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7440-38-2	Arsenic	2.48	ug/L	J	2	5	5	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7440-39-3	Barium	6.98	ug/L		1	5	5	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-41-7	Beryllium	1	ug/L	U	1	5	5	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-42-8	Boron	15	ug/L	U	15	50	50	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-43-9	Cadmium	0.30	ug/L	U	0.3	1	1	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7440-70-2	Calcium	7590	ug/L		50	200	200	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-47-3	Chromium	3	ug/L	U	3	10	10	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7440-48-4	Cobalt	1	ug/L	U	1	5	5	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-50-8	Copper	3	ug/L	U	3	10	10	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7439-89-6	Iron	30	ug/L	U	30	100	100	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7439-92-1	Lead	0.50	ug/L	U	0.5	2	2	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7439-95-4	Magnesium	2900	ug/L		110	300	300	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7439-96-5	Manganese	2	ug/L	U	2	10	10	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7439-98-7	Molybdenum	0.986	ug/L		0.2	0.5	0.5	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7440-02-0	Nickel	0.976	ug/L	J	0.6	2	2	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7440-09-7	Potassium	2190	ug/L		50	150	150	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7782-49-2	Selenium	2	ug/L	U	2	5	5	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7631-86-9	Silica	54500	ug/L		53	213	213	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-22-4	Silver	0.30	ug/L	U	0.3	1	1	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7440-23-5	Sodium	7990	ug/L		100	300	300	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-24-6	Strontium	43.1	ug/L		1	5	5	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-28-0	Thallium	0.60	ug/L	U	0.6	2	2	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7440-31-5	Tin	2.5	ug/L	U*	2.5	10	10	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-61-1	Uranium	0.282	ug/L		0.067	0.2	0.2	1	MS	BAJ	02/27/18 21:16	180227-2	1741864
7440-62-2	Vanadium	8.5	ug/L		1	5	5	1	P	TXT1	03/07/18 19:14	030718-1	1741857
7440-66-6	Zinc	8.94	ug/L	J	3.3	10	10	1	P	TXT1	03/07/18 19:14	030718-1	1741857

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811**CONTRACT:** ESHL00114**METHOD TYPE:****SAMPLE ID:** 444566011**BASIS:** As Received**DATE COLLECTED** 21-FEB-18**CLIENT ID:** CAWA-18-55**LEVEL:** Low**DATE RECEIVED** 23-FEB-18**MATRIX:** W**%SOLIDS:** 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
	Hardness as CaCO3	30.9	mg/L		0.453	1.24	1.24	1		TXT1	03/14/18 13:54		1747180

Prep Information:

Analytical Batch	Prep Batch	Prep Method	Initial wt./vol.	Units	Final wt./vol.	Units	Date	Analyst
1741857	1741856	SW846 3005A	50	mL	50	mL	02/23/18	JXM8
1741864	1741863	SW846 3005A	50	mL	50	mL	02/23/18	JXM8
1742549	1742548	EPA 245.1/245.2 Prep	20	mL	20	mL	02/27/18	AXS5

***Analytical Methods:**

P SW846 3005A/6010C
MS SW846 3005A/6020A
AV EPA 245.2 1974

METALS
-1-
INORGANICS ANALYSIS DATA PACKAGE

SDG No: 2018-1811**CONTRACT:** ESHL00114**METHOD TYPE:** EPA**SAMPLE ID:** 444566012**BASIS:** As Received**DATE COLLECTED** 21-FEB-18**CLIENT ID:** CAWA-18-56**LEVEL:** Low**DATE RECEIVED** 23-FEB-18**MATRIX:** W**%SOLIDS:** 0

CAS	Analyte	Result	Units	Qual	MDL	PQL	CRDL	DF	M*	Analyst	Run Date	Analytical Run	Analytical Batch
7439-97-6	Mercury	0.067	ug/L	U	0.067	0.2	0.2	1	AV	MTM1	02/28/18 10:38	022818W1-3	1742549

Prep Information:

Analytical Batch	Prep Batch	Prep Method	Initial wt./vol.	Units	Final wt./vol.	Units	Date	Analyst
1742549	1742548	EPA 245.1/245.2 Prep	20	mL	20	mL	02/27/18	AXS5

***Analytical Methods:**

AV EPA 245.2 1974

Quality Control Summary

METALS
-3b-
PREPARATION BLANK SUMMARY

SDG NO. 2018-1811
Contract: ESHL00114
Matrix: W

<u>Sample ID</u>	<u>Analyte</u>	<u>Result</u>	<u>Units</u>	<u>Acceptance Window</u>	<u>Conc Qual</u>	<u>M*</u>	<u>MDL</u>	<u>RDL</u>
1203978044	Aluminum	68	ug/L	+/-200	U	P	68	200
	Barium	1	ug/L	+/-5	U	P	1	5
	Beryllium	1	ug/L	+/-5	U	P	1	5
	Boron	15	ug/L	+/-50	U	P	15	50
	Calcium	50	ug/L	+/-200	U	P	50	200
	Cobalt	1	ug/L	+/-5	U	P	1	5
	Copper	3	ug/L	+/-10	U	P	3	10
	Iron	30	ug/L	+/-100	U	P	30	100
	Magnesium	110	ug/L	+/-300	U	P	110	300
	Manganese	2	ug/L	+/-10	U	P	2	10
	Potassium	50	ug/L	+/-150	U	P	50	150
	Silica	53	ug/L	+/-213	U	P	53	213
	Sodium	100	ug/L	+/-300	U	P	100	300
	Strontium	1	ug/L	+/-5	U	P	1	5
	Tin	2.5	ug/L	+/-10	U	P	2.5	10
	Vanadium	1	ug/L	+/-5	U	P	1	5
	Zinc	3.3	ug/L	+/-10	U	P	3.3	10
1203978054	Antimony	1	ug/L	+/-3	U	MS	1	3
	Arsenic	2	ug/L	+/-5	U	MS	2	5
	Cadmium	0.3	ug/L	+/-1	U	MS	0.3	1
	Chromium	3	ug/L	+/-10	U	MS	3	10
	Lead	0.5	ug/L	+/-2	U	MS	0.5	2
	Molybdenum	0.2	ug/L	+/-0.5	U	MS	0.2	0.5
	Nickel	0.6	ug/L	+/-2	U	MS	0.6	2
	Selenium	2	ug/L	+/-5	U	MS	2	5
	Silver	0.3	ug/L	+/-1	U	MS	0.3	1
	Thallium	0.6	ug/L	+/-2	U	MS	0.6	2
	Uranium	0.067	ug/L	+/-0.2	U	MS	0.067	0.2
1203979604	Mercury	0.067	ug/L	+/-0.2	U	AV	0.067	0.2

***Analytical Methods:**

P SW846 3005A/6010C
MS SW846 3005A/6020A
AV EPA 245.1/245.2

METALS

-5a-

Matrix Spike Summary

SDG NO. 2018-1811 Client ID: CAPA-18-1S

Contract: ESHL00114 Level: Low

Matrix: WATER % Solids:

Sample ID: 444566001 Spike ID: 1203978047

<u>Analyte</u>	<u>Units</u>	<u>Acceptance Limit</u>	<u>Spiked Result</u>	<u>C</u>	<u>Sample Result</u>	<u>C</u>	<u>Spike Added</u>	<u>% Recovery</u>	<u>Qual</u>	<u>M*</u>
Beryllium	ug/L	75-125	465		1	U	500	93.1		P
Boron	ug/L	75-125	463		15	U	500	92.6		P
Calcium	ug/L		26900		23000		5000	78.8	N/A	P
Cobalt	ug/L	75-125	457		1	U	500	91.4		P
Copper	ug/L	75-125	466		3	U	500	93.2		P
Iron	ug/L	75-125	4830		60	J	5000	95.5		P
Magnesium	ug/L	75-125	10700		6040		5000	93.6		P
Manganese	ug/L	75-125	458		2	U	500	91.5		P
Potassium	ug/L	75-125	7900		3330		5000	91.4		P
Silica	ug/L	75-125	48200		39800		10700	78.2		P
Sodium	ug/L		27800		24300		5000	69.7	N/A	P
Strontium	ug/L	75-125	582		131		500	90.3		P
Tin	ug/L	75-125	459		2.5	U	500	91.8		P
Vanadium	ug/L	75-125	467		2.54	J	500	92.9		P
Zinc	ug/L	75-125	449		3.3	U	500	89.2		P
Barium	ug/L	75-125	533		75.4		500	91.6		P
Aluminum	ug/L	75-125	4780		109	J	5000	93.5		P

*Analytical Methods:

P SW846 3005A/6010C

METALS

-5a-

Matrix Spike Summary

SDG NO. 2018-1811 Client ID: CAPA-18-1S

Contract: ESHL00114 Level: Low

Matrix: WATER % Solids:

Sample ID: 444566001 Spike ID: 1203978057

<u>Analyte</u>	<u>Units</u>	<u>Acceptance Limit</u>	<u>Spiked Result</u>	<u>C</u>	<u>Sample Result</u>	<u>C</u>	<u>Spike Added</u>	<u>% Recovery</u>	<u>Qual</u>	<u>M*</u>
Antimony	ug/L	75-125	48.5		1	U	50	96		MS
Arsenic	ug/L	75-125	51.3		2	U	50	99		MS
Cadmium	ug/L	75-125	49		0.3	U	50	98.1		MS
Chromium	ug/L	75-125	51.7		3	U	50	101		MS
Lead	ug/L	75-125	47.8		0.5	U	50	95.5		MS
Molybdenum	ug/L	75-125	50.8		1.14		50	99.4		MS
Nickel	ug/L	75-125	51.6		0.6	U	50	102		MS
Selenium	ug/L	75-125	51.4		2	U	50	101		MS
Silver	ug/L	75-125	49.9		0.3	U	50	99.9		MS
Thallium	ug/L	75-125	46.6		0.6	U	50	93.2		MS
Uranium	ug/L	75-125	47.5		0.494		50	94		MS

*Analytical Methods:

MS SW846 3005A/6020A

METALS

-5a-

Matrix Spike Summary

SDG NO. 2018-1811 **Client ID:** CAPA-18-1S**Contract:** ESHL00114 **Level:** Low**Matrix:** WATER **% Solids:****Sample ID:** 444566001 **Spike ID:** 1203979607

<u>Analyte</u>	<u>Units</u>	<u>Acceptance Limit</u>	<u>Spiked Result</u>	<u>C</u>	<u>Sample Result</u>	<u>C</u>	<u>Spike Added</u>	<u>% Recovery</u>	<u>Qual</u>	<u>M*</u>
Mercury	ug/L	75-125	2.01		0.067	U	2	100		AV

*Analytical Methods:

AV EPA 245.1/245.2

Metals
–6–
Duplicate Sample Summary

SDG No.: 2018–1811

Lab Code: GEL

Contract: ESHL00114

Client ID: CAPA–18–1D

Matrix: WATER

Level: Low

Sample ID: 444566001

Duplicate ID: 1203978046

Percent Solids for Dup: N/A

Analyte	Units	Acceptance Limit	Sample Result	C	Duplicate Result	C	RPD	Qual	M*
Aluminum	ug/L	+/-200	109 J		106 J		3.09		P
Barium	ug/L	+/-20%	75.4		74.6		1.05		P
Beryllium	ug/L		1 U		1 U				P
Boron	ug/L		15 U		15 U				P
Calcium	ug/L	+/-20%	23000		22800		.975		P
Cobalt	ug/L		1 U		1 U				P
Copper	ug/L		3 U		3 U				P
Iron	ug/L	+/-100	60 J		54.8 J		9.17		P
Magnesium	ug/L	+/-20%	6040		5950		1.45		P
Manganese	ug/L		2 U		2 U				P
Potassium	ug/L	+/-20%	3330		3320		.379		P
Silica	ug/L	+/-20%	39800		39100		1.89		P
Sodium	ug/L	+/-20%	24300		24200		.676		P
Strontium	ug/L	+/-20%	131		131		.366		P
Tin	ug/L	+/-10	-1.1 U		18.2		226	*	P
Vanadium	ug/L	+/-5	2.54 J		2.04 J		21.9		P
Zinc	ug/L		3.3 U		5.37 J		200		P

*Analytical Methods:

P SW846 3005A/6010C

Metals
-6-
Duplicate Sample Summary

SDG No.: 2018-1811

Lab Code: GEL

Contract: ESHL00114

Client ID: CAPA-18-1D

Matrix: WATER

Level: Low

Sample ID: 444566001

Duplicate ID: 1203978056

Percent Solids for Dup: N/A

Analyte	Units	Acceptance Limit	Sample Result	C	Duplicate Result	C	RPD	Qual	M*
Antimony	ug/L		1 U		1 U				MS
Arsenic	ug/L		2 U		2.02 J		200		MS
Cadmium	ug/L		0.3 U		0.3 U				MS
Chromium	ug/L		3 U		3 U				MS
Lead	ug/L		0.5 U		0.5 U				MS
Molybdenum	ug/L	+/- .5	1.14		1.12		2.57		MS
Nickel	ug/L		0.6 U		0.6 U				MS
Selenium	ug/L		2 U		2 U				MS
Silver	ug/L		0.3 U		0.3 U				MS
Thallium	ug/L		0.6 U		0.6 U				MS
Uranium	ug/L	+/- .2	0.494		0.498		.806		MS

*Analytical Methods:

MS SW846 3005A/6020A

Metals
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Duplicate Sample Summary

SDG No.: 2018–1811**Lab Code:** GEL**Contract:** ESHL00114**Client ID:** CAPA–18–1D**Matrix:** WATER**Level:** Low**Sample ID:** 444566001**Duplicate ID:** 1203979606**Percent Solids for Dup:** N/A

Analyte	Units	Acceptance Limit	Sample Result	C	Duplicate Result	C	RPD	Qual	M*
Mercury	ug/L		0.067	U	0.067	U			AV

*Analytical Methods:

AV EPA 245.1/245.2

METALS

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Laboratory Control Sample Summary

SDG NO. 2018-1811

Contract: ESHL00114

Aqueous LCS Source:OS2I

Solid LCS Source:

<u>Sample ID</u>	<u>Analyte</u>	<u>Units</u>	<u>True Value</u>	<u>Result</u>	<u>C</u>	<u>% Recovery</u>	<u>Acceptance Limit</u>	<u>M*</u>
1203978045								
	Aluminum	ug/L	5000	4800		96	80-120	P
	Barium	ug/L	500	477		95.4	80-120	P
	Beryllium	ug/L	500	478		95.6	80-120	P
	Boron	ug/L	500	472		94.4	80-120	P
	Calcium	ug/L	5000	4970		99.4	80-120	P
	Cobalt	ug/L	500	475		95	80-120	P
	Copper	ug/L	500	474		94.8	80-120	P
	Iron	ug/L	5000	4920		98.4	80-120	P
	Magnesium	ug/L	5000	5040		101	80-120	P
	Manganese	ug/L	500	481		96.2	80-120	P
	Potassium	ug/L	5000	4760		95.3	80-120	P
	Silica	ug/L	10700	9800		91.6	80-120	P
	Sodium	ug/L	5000	4680		93.6	80-120	P
	Strontium	ug/L	500	474		94.9	80-120	P
	Tin	ug/L	500	478		95.5	80-120	P
	Vanadium	ug/L	500	478		95.6	80-120	P
	Zinc	ug/L	500	461		92.3	80-120	P

*Analytical Methods:

P SW846 3005A/6010C

METALS

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Laboratory Control Sample Summary

SDG NO. 2018-1811

Contract: ESHL00114

Aqueous LCS Source: Inorganic Ventures

Solid LCS Source:

<u>Sample ID</u>	<u>Analyte</u>	<u>Units</u>	<u>True Value</u>	<u>Result</u>	<u>C</u>	<u>% Recovery</u>	<u>Acceptance Limit</u>	<u>M*</u>
1203978055								
	Cadmium	ug/L	50	48.6		97.3	80-120	MS
	Chromium	ug/L	50	49.5		99	80-120	MS
	Lead	ug/L	50	48.5		97.1	80-120	MS
	Molybdenum	ug/L	50	49.3		98.6	80-120	MS
	Nickel	ug/L	50	50.6		101	80-120	MS
	Selenium	ug/L	50	51.4		103	80-120	MS
	Silver	ug/L	50	49.8		99.6	80-120	MS
	Thallium	ug/L	50	47.7		95.5	80-120	MS
	Uranium	ug/L	50	46.8		93.6	80-120	MS
	Antimony	ug/L	50	48		96	80-120	MS
	Arsenic	ug/L	50	49.3		98.6	80-120	MS

*Analytical Methods:

MS SW846 3005A/6020A

METALS

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Laboratory Control Sample Summary

SDG NO. 2018-1811

Contract: ESHL00114

Aqueous LCS Source: GEL

Solid LCS Source:

<u>Sample ID</u>	<u>Analyte</u>	<u>Units</u>	<u>True Value</u>	<u>Result</u>	<u>C</u>	<u>% Recovery</u>	<u>Acceptance Limit</u>	<u>M*</u>
1203979605	Mercury	ug/L	2	2.13		107	85-115	AV

*Analytical Methods:

AV EPA 245.1/245.2

METALS

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Serial Dilution Sample Summary

SDG NO. 2018-1811

Client ID: CAPA-18-1L

Contract: ESHL00114

Matrix: LIQUID

Level: Low

Sample ID: 444566001

Serial Dilution ID: 1203978048

<u>Analyte</u>	<u>Initial Value ug/L</u>	<u>C</u>	<u>Serial Value ug/L</u>	<u>C</u>	<u>% Difference</u>	<u>Qual</u>	<u>Acceptance Limit</u>	<u>M*</u>
Aluminum	109	J	340	U	.931			P
Barium	75.4		76.4		1.389		10	P
Beryllium	1	U	5	U				P
Boron	15	U	75	U				P
Calcium	23000		22800		.905		10	P
Cobalt	1	U	5	U				P
Copper	3	U	15	U				P
Iron	60	J	150	U	12.081			P
Magnesium	6040		6070		.406		10	P
Manganese	2	U	10	U				P
Potassium	3330		3280		1.567		10	P
Silica	39800		39100		1.683		10	P
Sodium	24300		24200		.421		10	P
Strontium	131		131		.222		10	P
Tin	2.5	U	12.5	U				P
Vanadium	2.54	J	5	U	16.168			P
Zinc	3.3	U	16.5	U				P

*Analytical Methods:

P SW846 3005A/6010C

METALS

-9-

Serial Dilution Sample Summary

SDG NO. 2018-1811

Client ID: CAPA-18-1L

Contract: ESHL00114

Matrix: LIQUID

Level: Low

Sample ID: 444566001

Serial Dilution ID: 1203978058

<u>Analyte</u>	<u>Initial Value ug/L</u>	<u>C</u>	<u>Serial Value ug/L</u>	<u>C</u>	<u>% Difference</u>	<u>Qual</u>	<u>Acceptance Limit</u>	<u>M*</u>
Antimony	1	U	5	U				MS
Arsenic	2	U	10	U				MS
Cadmium	.3	U	1.5	U				MS
Chromium	3	U	15	U				MS
Lead	.5	U	2.5	U				MS
Molybdenum	1.14		1.37	J	19.755			MS
Nickel	.6	U	4.06	J				MS
Selenium	2	U	10	U				MS
Silver	.3	U	1.5	U				MS
Thallium	.6	U	3	U				MS
Uranium	.494		.51	J	3.239			MS

*Analytical Methods:

MS SW846 3005A/6020A

METALS

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Serial Dilution Sample Summary

SDG NO. 2018-1811 **Client ID:** CAPA-18-1L**Contract:** ESHL00114**Matrix:** LIQUID **Level:** Low**Sample ID:** 444566001 **Serial Dilution ID:** 1203979608

<u>Analyte</u>	<u>Initial Value ug/L</u>	<u>C</u>	<u>Serial Value ug/L</u>	<u>C</u>	<u>% Difference</u>	<u>Qual</u>	<u>Acceptance Limit</u>	<u>M*</u>
Mercury	.067	U	.335	U				AV

*Analytical Methods:

AV EPA 245.1/245.2

General Chem Analysis

Case Narrative

**General Chemistry
Technical Case Narrative
ARS International, LLC (ARSL)
SDG #: 2018-1811
Work Order #: 444566**

Method/Analysis Information

Product: Carbon and Total Organic

Analytical Batch: 1741854

Method: SW 9060 Total Organic Carbon

Sample Analysis

The following samples were analyzed using the analytical protocol as established in SW-846:9060:

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566012	CAWA-18-56
1203980262	Method Blank (MB)
1203980263	Laboratory Control Sample (LCS)
1203980265	444446006(CAMO-18-28) Sample Duplicate (DUP)
1203980269	444446006(CAMO-18-28) Post Spike (PS)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-093 REV# 15.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC, and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Carbon analysis was performed on a O-I Analytical 1030W Carbon Analyzer.

Initial Calibration

All initial calibration requirements have been met for this SDG.

Continuing Calibration Blanks

All continuing calibration blanks (CCBs) associated with reported data from this batch were within acceptance limits.

Calibration Verification Information (CCV)

All continuing calibration verification standards (CCVs) associated with reported data from this batch were within acceptance limits.

Quality Control (QC) Information**Method Blank (MB) Statement**

The MB analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Sample 444446006 (CAMO-18-28) was selected for QC analysis.

Matrix Spike (MS)/Post Spike (PS) Recovery Statement

The MS/PS recovery for this sample set was within the required acceptance limits where applicable.

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Preservation/Integrity

All the samples from this sample group met the preservation and integrity requirements of the method.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information**Additional Comments**

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Method/Analysis Information

Product:	Cyanide and Total		
Analytical Batch:	1740518	Method:	WSP-CN(T)
Prep Batch :	1740517	Method:	EPA 335.4

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA 335.4 1993:

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566012	CAWA-18-56
1203975195	Method Blank (MB)
1203975196	Laboratory Control Sample (LCS)
1203975197	444396002(CAWA-18-28) Sample Duplicate (DUP)
1203975198	444396002(CAWA-18-28) Matrix Spike (MS)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-095 REV# 21.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Flow Injection analysis was performed on a Lachat QuickChem FIA+ 8000 Series.

Initial Calibration

All initial calibration requirements have been met for this SDG.

Continuing Calibration Blanks

All continuing calibration blanks (CCBs) associated with reported data from this batch were within acceptance limits.

Calibration Verification Information (CCV)

All continuing calibration verification standards (CCVs) associated with reported data from this batch were within acceptance limits.

Y Intercept Rule

The absolute value of the intercept is less than 3 times the MDL.

Quality Control (QC) Information**Method Blank (MB) Statement**

The MB analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Sample 444396002 (CAWA-18-28) was selected for QC analysis.

Matrix Spike (MS)/Post Spike (PS) Recovery Statement

The MS/PS recovery for this sample set was within the required acceptance limits where applicable.

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Preservation/Integrity

All the samples from this sample group met the preservation and integrity requirements of the method.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information**Additional Comments**

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Method/Analysis Information

Product: Ion Chromatography

Analytical Batch: 1741949

Method: WSP-ANIONS

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA:300.0:

Sample ID	Client ID
444566001	CAPA-18-1
444566005	CAPA-18-10
444566011	CAWA-18-55
1203978236	Method Blank (MB)
1203978237	Laboratory Control Sample (LCS)
1203978238	444565001(CTUA-17-151327) Sample Duplicate (DUP)
1203978239	444565001(CTUA-17-151327) Post Spike (PS)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-086 REV# 25.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Ion Chromatography analysis was performed on a Dionex ICS-3000 Ion Chromatograph.

Initial Calibration

All initial calibration requirements have been met for this SDG.

Continuing Calibration Blanks

All continuing calibration blanks (CCBs) associated with reported data from this batch were within acceptance limits.

Calibration Verification Information (CCV)

All continuing calibration verification standards (CCVs) associated with reported data from this batch were within acceptance limits.

Y Intercept Rule

The absolute value of the intercept is less than 3 times the MDL.

Quality Control (QC) Information

Method Blank (MB) Statement

The MB analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Sample 444565001 (CTUA-17-151327) was selected for QC analysis.

Matrix Spike (MS)/Post Spike (PS) Recovery Statement

The percent recoveries (%R) obtained from the spike analyses are evaluated when the sample concentration is less than four times (4X) the spike concentration added. The matrix spike recovered outside of the established acceptance limits due to matrix interference and/or non-homogeneity.

Analyte	Sample	Value
Sulfate	1203978239 (CTUA-17-151327PS)	130* (75%-125%)

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Dilutions

The following samples 1203978238 (CTUA-17-151327DUP), 1203978239 (CTUA-17-151327PS), 444566001 (CAPA-18-1) and 444566005 (CAPA-18-10) were diluted because target analyte concentrations exceeded the calibration range. Dilutions may be required for many reasons, including to minimize matrix interferences or to bring over range target analyte concentrations into the linear calibration range.

Analyte	444566	
	001	005
Chloride	5X	5X

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information

Manual Integrations

Samples 1203978238 (CTUA-17-151327DUP), 1203978239 (CTUA-17-151327PS), 444566001 (CAPA-18-1), 444566005 (CAPA-18-10) and 444566011 (CAWA-18-55) were manually integrated to correctly position the baseline as set in the calibration standards.

Additional Comments

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Method/Analysis Information

Product: Ammonia Nitrogen
Analytical Batch: 1741276 **Method:** NH3
Prep Batch : 1741272 **Method:** EPA 350.1 Prep

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA:350.1:

Sample ID	Client ID
444566001	CAPA-18-1
444566005	CAPA-18-10
444566011	CAWA-18-55
1203976869	Method Blank (MB)
1203976870	Laboratory Control Sample (LCS)
1203976872	444566001(CAPA-18-1) Sample Duplicate (DUP)
1203976874	444566001(CAPA-18-1) Matrix Spike (MS)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-106 REV# 10.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Nutrient analysis was performed on a Lachat QuickChem FIA+ 8000 Series.

Initial Calibration

All initial calibration requirements have been met for this SDG.

Calibration Verification Information

All associated calibration verification standard(s) (ICV or CCV) met the acceptance criteria.

Continuing Calibration Blanks

All continuing calibration blanks (CCBs) associated with reported data from this batch were within acceptance limits.

Calibration Verification Information (CCV)

All continuing calibration verification standards (CCVs) associated with reported data from this batch were within acceptance limits.

Y Intercept Rule

The absolute value of the intercept is less than 3 times the MDL.

Quality Control (QC) Information**Method Blank (MB) Statement**

The MB analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Sample 444566001 (CAPA-18-1) was selected for QC analysis.

Matrix Spike (MS)/Post Spike (PS) Recovery Statement

The MS/PS recovery for this sample set was within the required acceptance limits where applicable.

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Preservation/Integrity

All the samples from this sample group met the preservation and integrity requirements of the method.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information**Additional Comments**

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Method/Analysis Information

Product:	Total Kjeldahl Nitrogen		
Analytical Batch:	1742074	Method:	TKN
Prep Batch :	1742073	Method:	EPA 351.2 Prep

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA:351.2:

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566012	CAWA-18-56
1203978464	Method Blank (MB)
1203978465	Laboratory Control Sample (LCS)
1203978466	444396002(CAWA-18-28) Sample Duplicate (DUP)
1203978468	444396006(CAWA-18-36) Sample Duplicate (DUP)
1203978467	444396002(CAWA-18-28) Matrix Spike (MS)
1203978469	444396006(CAWA-18-36) Matrix Spike (MS)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-104 REV# 15.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Nutrient analysis was performed on a Lachat QuickChem FIA+ 8000 Series.

Initial Calibration

All initial calibration requirements have been met for this SDG.

Calibration Verification Information

All associated calibration verification standard(s) (ICV or CCV) met the acceptance criteria.

Continuing Calibration Blanks

All continuing calibration blanks (CCBs) associated with reported data from this batch were within acceptance limits.

Calibration Verification Information (CCV)

All continuing calibration verification standards (CCVs) associated with reported data from this batch were within acceptance limits.

Y Intercept Rule

The absolute value of the intercept is less than 3 times the MDL.

Quality Control (QC) Information**Method Blank (MB) Statement**

The MB analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Samples 444396002 (CAWA-18-28) and 444396006 (CAWA-18-36) were selected for QC analysis.

Matrix Spike (MS)/Post Spike (PS) Recovery Statement

The percent recoveries (%R) obtained from the spike analyses are evaluated when the sample concentration is less than four times (4X) the spike concentration added. The matrix spike recovered outside of the established acceptance limits due to matrix interference and/or non-homogeneity.

Analyte	Sample	Value
Nitrogen, Total Kjeldahl	1203978467 (CAWA-18-28MS)	50.1* (90%-110%)
	1203978469 (CAWA-18-36MS)	75.1* (90%-110%)

Duplicate Relative Percent Difference (RPD) Statement

The Relative Percent Difference (RPD) between the sample and duplicate falls outside of the established acceptance limits because of the heterogeneous matrix of the sample:

Analyte	Sample	Value
Nitrogen, Total Kjeldahl	1203978466 (CAWA-18-28DUP)	abs(-.0202 - .164)* (+/- .1 mg/L)

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Preservation/Integrity

All the samples from this sample group met the preservation and integrity requirements of the method.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

Samples 1203978467 (CAWA-18-28MS), 1203978468 (CAWA-18-36DUP) and 1203978469 (CAWA-18-36MS) were re-analyzed due to CCV failure. The reanalysis data with passing instrument QC was reported.

Miscellaneous Information**Additional Comments**

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

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Method/Analysis Information

Product: Nitrate Nitrite by Cadmium Reduction

Analytical Batch: 1742109

Method: NO3NO2

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA:353.2:

Sample ID	Client ID
444566001	CAPA-18-1
444566005	CAPA-18-10
444566011	CAWA-18-55
1203978566	Method Blank (MB)
1203978567	Laboratory Control Sample (LCS)
1203978569	444396001(CAWA-18-27) Sample Duplicate (DUP)
1203978571	444396001(CAWA-18-27) Post Spike (PS)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-128 REV# 10.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Nutrient analysis was performed on a Lachat QuickChem FIA+ 8500 Series.

Initial Calibration

All initial calibration requirements have been met for this SDG.

Calibration Verification Information

All associated calibration verification standard(s) (ICV or CCV) met the acceptance criteria.

Continuing Calibration Blanks

All continuing calibration blanks (CCBs) associated with reported data from this batch were within acceptance limits.

Calibration Verification Information (CCV)

All continuing calibration verification standards (CCVs) associated with reported data from this batch were within

acceptance limits.

Y Intercept Rule

The absolute value of the intercept is less than 3 times the MDL.

Quality Control (QC) Information

Method Blank (MB) Statement

The MB analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Sample 444396001 (CAWA-18-27) was selected for QC analysis.

Matrix Spike (MS)/Post Spike (PS) Recovery Statement

The MS/PS recoveries for this sample set were within the required acceptance limits where applicable.

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Preservation/Integrity

All the samples from this sample group met the preservation and integrity requirements of the method.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information

Additional Comments

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages

electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Method/Analysis Information

Product:	Total Phosphorus		
Analytical Batch:	1742097	Method:	PO4
Prep Batch :	1742096	Method:	EPA 365.4 Prep

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA 365.4 1974:

Sample ID	Client ID
444566001	CAPA-18-1
444566005	CAPA-18-10
444566011	CAWA-18-55
1203978529	Method Blank (MB)
1203978530	Laboratory Control Sample (LCS)
1203978531	444396001(CAWA-18-27) Sample Duplicate (DUP)
1203978532	444396001(CAWA-18-27) Matrix Spike (MS)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-103 REV# 11.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Nutrient analysis was performed on a Lachat QuickChem FIA+ 8000 Series.

Initial Calibration

All initial calibration requirements have been met for this SDG.

Continuing Calibration Blanks

All continuing calibration blanks (CCBs) associated with reported data from this batch were within acceptance limits.

Calibration Verification Information (CCV)

All continuing calibration verification standards (CCVs) associated with reported data from this batch were within acceptance limits.

Y Intercept Rule

The absolute value of the intercept is less than 3 times the MDL.

Quality Control (QC) Information**Method Blank (MB) Statement**

The MB analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Sample 444396001 (CAWA-18-27) was selected for QC analysis.

Matrix Spike (MS)/Post Spike (PS) Recovery Statement

The MS/PS recovery for this sample set was within the required acceptance limits where applicable.

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Preservation/Integrity

All the samples from this sample group met the preservation and integrity requirements of the method.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information**Additional Comments**

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Method/Analysis Information

Product: Solids and Total Dissolved

Analytical Batch: 1742467

Method: TDS

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA:160.1:

Sample ID	Client ID
444566001	CAPA-18-1
444566005	CAPA-18-10
444566011	CAWA-18-55
1203979421	Method Blank (MB)
1203979422	Laboratory Control Sample (LCS)
1203979423	444446003(CAMO-18-19) Sample Duplicate (DUP)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-001 REV# 15.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Solids analysis was performed on a Sartorius Balance BAL216. Solids lab

Initial Calibration

All initial calibration requirements have been met for this SDG.

Quality Control (QC) Information

Method Blank (MB) Statement

The MB analyzed with this SDG met the acceptance criteria.

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Consecutive Weight Checks

All consecutive weight checks were met.

Quality Control (QC) Designation

Sample 444446003 (CAMO-18-19) was selected for QC analysis.

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information**Additional Comments**

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Method/Analysis Information

Product: **Specific Conductivity**

Analytical Batch: 1742552 **Method:** EPA120.1 Specific Conductivity

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA:120.1:

Sample ID	Client ID
444566001	CAPA-18-1
444566005	CAPA-18-10
444566011	CAWA-18-55
1203979622	Laboratory Control Sample (LCS)
1203979623	444446001(CAMO-18-16) Sample Duplicate (DUP)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-009 REV# 16.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Titration and Ion analysis was performed on a Thermo Scientific Orion Star A212 Conductivity Meter.

Initial Calibration

All initial calibration requirements have been met for this SDG.

Initial Standardization

The titrant was properly standardized

Quality Control (QC) Information

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Sample 444446001 (CAMO-18-16) was selected for QC analysis.

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information**Additional Comments**

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Method/Analysis Information

Product: pH
Analytical Batch: 1741049 **Method:** PH

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA 150.1 1982:

Sample ID	Client ID
444566001	CAPA-18-1
444566005	CAPA-18-10
444566011	CAWA-18-55
1203976494	Laboratory Control Sample (LCS)
1203976495	444026001(CAMO-18-1) Sample Duplicate (DUP)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-008 REV# 23.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Titration and Ion analysis was performed on a Thermo Orion Star A111. Immediates

Initial Calibration

All initial calibration requirements have been met for this SDG.

Initial Standardization

The titrant was properly standardized

Quality Control (QC) Information

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Sample 444026001 (CAMO-18-1) was selected for QC analysis.

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

Samples (See Below) were received by the laboratory outside of the method specified holding time. The data is qualified.

Sample	Analyte	Value
1203976495 (CAMO-18-1DUP)	pH	Received 19-FEB-18, out of holding 14-FEB-18
444566001 (CAPA-18-1)	pH	Received 23-FEB-18, out of holding 21-FEB-18
444566005 (CAPA-18-10)	pH	Received 23-FEB-18, out of holding 21-FEB-18
444566011 (CAWA-18-55)	pH	Received 23-FEB-18, out of holding 21-FEB-18

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information**Additional Comments**

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Method/Analysis Information

Product: Alkalinity

Analytical Batch: 1741047 **Method:** EPA 310.1 Total Alkalinity

Sample Analysis

The following samples were analyzed using the analytical protocol as established in EPA:310.1:

Sample ID	Client ID
444566001	CAPA-18-1
444566005	CAPA-18-10
444566011	CAWA-18-55
1203976485	Laboratory Control Sample (LCS)
1203976486	444026001(CAMO-18-1) Sample Duplicate (DUP)
1203976487	444396029(CAWA-18-29) Sample Duplicate (DUP)
1203976488	444026001(CAMO-18-1) Matrix Spike (MS)
1203976489	444396029(CAWA-18-29) Matrix Spike (MS)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-GC-E-033 REV# 13.

Preparation/Analytical Method Verification

The SOP stated above has been prepared based on technical research and testing conducted by GEL Laboratories, LLC. and with guidance from the regulatory documents listed in this "Method/Analysis Information" section.

Calibration Information

The Titration and Ion analysis was performed on a Electronic bottle-top buret.

Initial Calibration

All initial calibration requirements have been met for this SDG.

Initial Standardization

The titrant was properly standardized

Quality Control (QC) Information

Laboratory Control Sample Duplicate (LCSD)

An LCSD was not used in place of matrix QC.

Laboratory Control Sample (LCS) Recovery

The LCS spike recovery met the acceptance limits.

Quality Control (QC) Designation

Samples 444026001 (CAMO-18-1) and 444396029 (CAWA-18-29) were selected for QC analysis.

Matrix Spike (MS)/Post Spike (PS) Recovery Statement

The MS/PS recoveries for this sample set were within the required acceptance limits where applicable.

Duplicate Relative Percent Difference (RPD) Statement

The RPD between the sample and its duplicate met the acceptance limits.

Technical Information

GEL assigns holding times based on the date and time of sample collection. Those holding times expressed in hours are calculated in the AlphaLims system by hours. Those holding times expressed as days expire at midnight on the day of expiration.

Holding Times

All samples in this SDG met the specified holding time.

Sample Dilutions

The samples in this SDG did not require dilutions.

Sample Re-analysis

The samples in this SDG did not require re-analysis.

Miscellaneous Information**Additional Comments**

Additional comments were not required for this SDG.

Electronic Packaging Comment

This data package was generated using an electronic data processing program referred to as virtual packaging. In an effort to increase quality and efficiency, the laboratory has developed systems to generate all data packages electronically. The following change from traditional packages should be noted:

Analyst/peer reviewer initials and dates are not present on the electronic data files. Presently, all initials and dates are present on the original raw data. These hard copies are temporarily stored in the laboratory. The data validator will always sign and date the case narrative. Data that are not generated electronically, such as hand written pages, will be scanned and inserted into the electronic package.

Certification Statement

Where the analytical method has been performed under NELAP certification, the analysis has met all of the requirements of the NELAC standard unless otherwise noted in the analytical case narrative.

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Qualifier Definition Report for

ARSL004 ARS International, LLC (ARS-LANS-MTOA6-25093-GEL)

Client SDG: 2018-1811 GEL Work Order: 444566

The Qualifiers in this report are defined as follows:

- * A quality control analyte recovery is outside of specified acceptance criteria
- H Analytical holding time was exceeded
- J Value is estimated
- U Analyte was analyzed for, but not detected above the MDL, MDA, MDC or LOD.

Review/Validation

GEL requires all analytical data to be verified by a qualified data reviewer. In addition, all CLP-like deliverables receive a third level review of the fractional data package.

The following data validator verified the information presented in this data report:

Signature:



Name: Kristen Mizzell

Date: 19 MAR 2018

Title: Team Leader

Sample Data Summary

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Certificate of Analysis

Report Date: March 19, 2018

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545
Contact: Ms. Nita Patel
Project: LANL- WQH Water Samples

Client SDG: 2018-1811

Client Sample ID: CAPA-18-1
Sample ID: 444566001
Matrix: W
Collect Date: 21-FEB-18 11:47
Receive Date: 23-FEB-18
Collector: Client

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	DL	RL	Units	PF	DF	Analyst	Date	Time	Batch	Method
Ion Chromatography												
WSP-ANIONS "As Received"												
Bromide	J	0.108	0.067	0.200	mg/L		1	JXH5	02/23/18	2023	1741949	1
Fluoride		0.231	0.033	0.100	mg/L		1					
Sulfate		17.1	0.133	0.400	mg/L		1					
Chloride		27.5	0.335	1.00	mg/L		5	JXH5	02/26/18	1246	1741949	2
Nutrient Analysis												
NH3 "As Received"												
Nitrogen, Ammonia		0.0759	0.017	0.050	mg/L	1.00	1	KLP1	02/27/18	1032	1741276	3
NO3NO2 "As Received"												
Nitrogen, Nitrate/Nitrite		1.40	0.017	0.050	mg/L		1	AXH3	02/26/18	0921	1742109	4
PO4 "As Received"												
Phosphorus, Total as P		0.0827	0.020	0.050	mg/L	1.00	1	KLP1	03/05/18	1225	1742097	5
Solids Analysis												
TDS "As Received"												
Total Dissolved Solids		194	3.40	14.3	mg/L			KLP1	02/27/18	1256	1742467	6
Titration and Ion Analysis												
EPA 310.1 Total Alkalinity "As Received"												
Alkalinity, Total as CaCO3		80.8	1.45	4.00	mg/L			RXB5	02/24/18	1515	1741047	7
Carbonate alkalinity (CaCO3)	U	ND	1.45	4.00	mg/L							
EPA120.1 Specific Conductivity "As Received"												
Conductivity		304	1.00	1.00	umhos/cm		1	HXC1	02/27/18	1317	1742552	8
PH "As Received"												
pH at Temp 16.5C	H	7.97	0.010	0.100	SU		1	RXB5	02/24/18	1517	1741049	9

The following Prep Methods were performed:

Method	Description	Analyst	Date	Time	Prep Batch
EPA 350.1 Prep	EPA 350.1 Ammonia Nitrogen Prep	AXH3	02/27/18	0702	1741272
EPA 365.4 Prep	EPA 365.4 Phosphorus, Total in liquid PR	AXH3	03/05/18	0900	1742096

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Certificate of Analysis

Report Date: March 19, 2018

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545
Contact: Ms. Nita Patel
Project: LANL- WQH Water Samples

Client SDG: 2018-1811

Client Sample ID: CAPA-18-1
Sample ID: 444566001

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	DL	RL	Units	PF	DF	Analyst	Date	Time	Batch	Method
The following Analytical Methods were performed:												
Method	Description					Analyst Comments						
1	EPA:300.0											
2	EPA:300.0											
3	EPA:350.1											
4	EPA:353.2											
5	EPA 365.4 1974											
6	EPA:160.1											
7	EPA:310.1											
8	EPA:120.1											
9	EPA 150.1 1982											

Notes:

Column headers are defined as follows:

DF: Dilution Factor

DL: Detection Limit

MDA: Minimum Detectable Activity

MDC: Minimum Detectable Concentration

Lc/LC: Critical Level

PF: Prep Factor

RL: Reporting Limit

SQL: Sample Quantitation Limit

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Certificate of Analysis

Report Date: March 19, 2018

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545
Contact: Ms. Nita Patel
Project: LANL- WQH Water Samples

Client SDG: 2018-1811

Client Sample ID: CAPA-18-2
Sample ID: 444566002
Matrix: W
Collect Date: 21-FEB-18 11:47
Receive Date: 23-FEB-18
Collector: Client

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	DL	RL	Units	PF	DF	Analyst	Date	Time	Batch	Method
Carbon Analysis												
SW 9060 Total Organic Carbon "As Received"												
Total Organic Carbon Average		1.69	0.330	1.00	mg/L		1	TSM	03/01/18	0008	1741854	1
Flow Injection Analysis												
WSP-CN(T) "As Received"												
Cyanide, Total	U	ND	1.67	5.00	ug/L	1.00	1	AXH3	02/28/18	0810	1740518	2
Nutrient Analysis												
TKN "As Received"												
Nitrogen, Total Kjeldahl	J	0.0408	0.033	0.100	mg/L	1.00	1	KLP1	02/28/18	1504	1742074	3

The following Prep Methods were performed:

Method	Description	Analyst	Date	Time	Prep Batch
EPA 335.4	EPA 335.4 Total Cyanide	AXH3	02/28/18	0630	1740517
EPA 351.2 Prep	EPA 351.2 Total Kjeldahl Nitrogen Prep	AXH3	02/27/18	1000	1742073

The following Analytical Methods were performed:

Method	Description	Analyst Comments
1	SW-846:9060	
2	EPA 335.4 1993	
3	EPA:351.2	

Notes:

Column headers are defined as follows:

DF: Dilution Factor	Lc/LC: Critical Level
DL: Detection Limit	PF: Prep Factor
MDA: Minimum Detectable Activity	RL: Reporting Limit
MDC: Minimum Detectable Concentration	SQL: Sample Quantitation Limit

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Certificate of Analysis

Report Date: March 19, 2018

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545
Contact: Ms. Nita Patel
Project: LANL- WQH Water Samples

Client SDG: 2018-1811

Client Sample ID: CAPA-18-10
Sample ID: 444566005
Matrix: W
Collect Date: 21-FEB-18 11:47
Receive Date: 23-FEB-18
Collector: Client

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	DL	RL	Units	PF	DF	Analyst	Date	Time	Batch	Method
Ion Chromatography												
WSP-ANIONS "As Received"												
Bromide	J	0.112	0.067	0.200	mg/L		1	JXH5	02/23/18	2054	1741949	1
Fluoride		0.269	0.033	0.100	mg/L		1					
Sulfate		17.1	0.133	0.400	mg/L		1					
Chloride		27.1	0.335	1.00	mg/L		5	JXH5	02/26/18	1317	1741949	2
Nutrient Analysis												
NH3 "As Received"												
Nitrogen, Ammonia	J	0.0284	0.017	0.050	mg/L	1.00	1	KLP1	02/27/18	1034	1741276	3
NO3NO2 "As Received"												
Nitrogen, Nitrate/Nitrite		1.40	0.017	0.050	mg/L		1	AXH3	02/26/18	0922	1742109	4
PO4 "As Received"												
Phosphorus, Total as P		0.0852	0.020	0.050	mg/L	1.00	1	KLP1	03/05/18	1226	1742097	5
Solids Analysis												
TDS "As Received"												
Total Dissolved Solids		186	3.40	14.3	mg/L			KLP1	02/27/18	1256	1742467	6
Titration and Ion Analysis												
EPA 310.1 Total Alkalinity "As Received"												
Alkalinity, Total as CaCO3		82.4	1.45	4.00	mg/L			RXB5	02/24/18	1520	1741047	7
Carbonate alkalinity (CaCO3)	U	ND	1.45	4.00	mg/L							
EPA120.1 Specific Conductivity "As Received"												
Conductivity		300	1.00	1.00	umhos/cm		1	HXC1	02/27/18	1318	1742552	8
PH "As Received"												
pH at Temp 16.3C	H	7.98	0.010	0.100	SU		1	RXB5	02/24/18	1517	1741049	9

The following Prep Methods were performed:

Method	Description	Analyst	Date	Time	Prep Batch
EPA 350.1 Prep	EPA 350.1 Ammonia Nitrogen Prep	AXH3	02/27/18	0702	1741272
EPA 365.4 Prep	EPA 365.4 Phosphorus, Total in liquid PR	AXH3	03/05/18	0900	1742096

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Certificate of Analysis

Report Date: March 19, 2018

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545
Contact: Ms. Nita Patel
Project: LANL- WQH Water Samples

Client SDG: 2018-1811

Client Sample ID: CAPA-18-10
Sample ID: 444566005

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	DL	RL	Units	PF	DF	Analyst	Date	Time	Batch	Method
The following Analytical Methods were performed:												
Method	Description					Analyst Comments						
1	EPA:300.0											
2	EPA:300.0											
3	EPA:350.1											
4	EPA:353.2											
5	EPA 365.4 1974											
6	EPA:160.1											
7	EPA:310.1											
8	EPA:120.1											
9	EPA 150.1 1982											

Notes:

Column headers are defined as follows:

DF: Dilution Factor

DL: Detection Limit

MDA: Minimum Detectable Activity

MDC: Minimum Detectable Concentration

Lc/LC: Critical Level

PF: Prep Factor

RL: Reporting Limit

SQL: Sample Quantitation Limit

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Certificate of Analysis

Report Date: March 19, 2018

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545
Contact: Ms. Nita Patel
Project: LANL- WQH Water Samples

Client SDG: 2018-1811

Client Sample ID: CAPA-18-12
Sample ID: 444566006
Matrix: W
Collect Date: 21-FEB-18 11:47
Receive Date: 23-FEB-18
Collector: Client

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	DL	RL	Units	PF	DF	Analyst	Date	Time	Batch	Method
Carbon Analysis												
SW 9060 Total Organic Carbon "As Received"												
Total Organic Carbon Average		1.68	0.330	1.00	mg/L		1	TSM	03/01/18	0048	1741854	1
Flow Injection Analysis												
WSP-CN(T) "As Received"												
Cyanide, Total	U	ND	1.67	5.00	ug/L	1.00	1	AXH3	02/28/18	0811	1740518	2
Nutrient Analysis												
TKN "As Received"												
Nitrogen, Total Kjeldahl	J	0.038	0.033	0.100	mg/L	1.00	1	KLP1	02/28/18	1505	1742074	3

The following Prep Methods were performed:

Method	Description	Analyst	Date	Time	Prep Batch
EPA 335.4	EPA 335.4 Total Cyanide	AXH3	02/28/18	0630	1740517
EPA 351.2 Prep	EPA 351.2 Total Kjeldahl Nitrogen Prep	AXH3	02/27/18	1000	1742073

The following Analytical Methods were performed:

Method	Description	Analyst Comments
1	SW-846:9060	
2	EPA 335.4 1993	
3	EPA:351.2	

Notes:

Column headers are defined as follows:

DF: Dilution Factor	Lc/LC: Critical Level
DL: Detection Limit	PF: Prep Factor
MDA: Minimum Detectable Activity	RL: Reporting Limit
MDC: Minimum Detectable Concentration	SQL: Sample Quantitation Limit

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Certificate of Analysis

Report Date: March 19, 2018

Company : Los Alamos National Laboratory
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Los Alamos, New Mexico 87545
Contact: Ms. Nita Patel
Project: LANL- WQH Water Samples

Client SDG: 2018-1811

Client Sample ID: CAWA-18-55
Sample ID: 444566011
Matrix: W
Collect Date: 21-FEB-18 10:44
Receive Date: 23-FEB-18
Collector: Client

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	DL	RL	Units	PF	DF	Analyst	Date	Time	Batch	Method
Ion Chromatography												
WSP-ANIONS "As Received"												
Bromide	U	ND	0.067	0.200	mg/L		1	JXH5	02/23/18	2124	1741949	1
Chloride		1.19	0.067	0.200	mg/L		1					
Fluoride		0.114	0.033	0.100	mg/L		1					
Sulfate		1.19	0.133	0.400	mg/L		1					
Nutrient Analysis												
NH3 "As Received"												
Nitrogen, Ammonia		0.0625	0.017	0.050	mg/L	1.00	1	KLP1	02/27/18	1035	1741276	2
NO3NO2 "As Received"												
Nitrogen, Nitrate/Nitrite		0.374	0.017	0.050	mg/L		1	AXH3	02/26/18	0923	1742109	3
PO4 "As Received"												
Phosphorus, Total as P		0.0814	0.020	0.050	mg/L	1.00	1	KLP1	03/05/18	1227	1742097	4
Solids Analysis												
TDS "As Received"												
Total Dissolved Solids		116	3.40	14.3	mg/L			KLP1	02/27/18	1256	1742467	5
Titration and Ion Analysis												
EPA 310.1 Total Alkalinity "As Received"												
Alkalinity, Total as CaCO3		48.6	1.45	4.00	mg/L			RXB5	02/24/18	1523	1741047	6
Carbonate alkalinity (CaCO3)	U	ND	1.45	4.00	mg/L							
EPA120.1 Specific Conductivity "As Received"												
Conductivity		101	1.00	1.00	umhos/cm		1	HXC1	02/27/18	1319	1742552	7
PH "As Received"												
pH at Temp 16.1C	H	7.70	0.010	0.100	SU		1	RXB5	02/24/18	1521	1741049	8

The following Prep Methods were performed:

Method	Description	Analyst	Date	Time	Prep Batch
EPA 350.1 Prep	EPA 350.1 Ammonia Nitrogen Prep	AXH3	02/27/18	0702	1741272
EPA 365.4 Prep	EPA 365.4 Phosphorus, Total in liquid PR	AXH3	03/05/18	0900	1742096

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Certificate of Analysis

Report Date: March 19, 2018

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545
Contact: Ms. Nita Patel
Project: LANL- WQH Water Samples

Client SDG: 2018-1811

Client Sample ID: CAWA-18-55
Sample ID: 444566011

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	DL	RL	Units	PF	DF	Analyst	Date	Time	Batch	Method
The following Analytical Methods were performed:												
Method	Description		Analyst Comments									
1	EPA:300.0											
2	EPA:350.1											
3	EPA:353.2											
4	EPA 365.4 1974											
5	EPA:160.1											
6	EPA:310.1											
7	EPA:120.1											
8	EPA 150.1 1982											

Notes:

Column headers are defined as follows:

DF: Dilution Factor	Lc/LC: Critical Level
DL: Detection Limit	PF: Prep Factor
MDA: Minimum Detectable Activity	RL: Reporting Limit
MDC: Minimum Detectable Concentration	SQL: Sample Quantitation Limit

GEL LABORATORIES LLC

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Certificate of Analysis

Report Date: March 19, 2018

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545

Contact: Ms. Nita Patel
Project: LANL- WQH Water Samples

Client SDG: 2018-1811

Client Sample ID: CAWA-18-56
Sample ID: 444566012
Matrix: W
Collect Date: 21-FEB-18 10:44
Receive Date: 23-FEB-18
Collector: Client

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	DL	RL	Units	PF	DF	Analyst	Date	Time	Batch	Method
Carbon Analysis												
SW 9060 Total Organic Carbon "As Received"												
Total Organic Carbon Average	J	0.339	0.330	1.00	mg/L		1	TSM	03/01/18	0128	1741854	1
Flow Injection Analysis												
WSP-CN(T) "As Received"												
Cyanide, Total	U	ND	1.67	5.00	ug/L	1.00	1	AXH3	02/28/18	0812	1740518	2
Nutrient Analysis												
TKN "As Received"												
Nitrogen, Total Kjeldahl	J	0.0474	0.033	0.100	mg/L	1.00	1	KLP1	02/28/18	1506	1742074	3

The following Prep Methods were performed:

Method	Description	Analyst	Date	Time	Prep Batch
EPA 335.4	EPA 335.4 Total Cyanide	AXH3	02/28/18	0630	1740517
EPA 351.2 Prep	EPA 351.2 Total Kjeldahl Nitrogen Prep	AXH3	02/27/18	1000	1742073

The following Analytical Methods were performed:

Method	Description	Analyst Comments
1	SW-846:9060	
2	EPA 335.4 1993	
3	EPA:351.2	

Notes:

Column headers are defined as follows:

DF: Dilution Factor	Lc/LC: Critical Level
DL: Detection Limit	PF: Prep Factor
MDA: Minimum Detectable Activity	RL: Reporting Limit
MDC: Minimum Detectable Concentration	SQL: Sample Quantitation Limit

Quality Control Summary

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QC Summary

Report Date: March 19, 2018

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Los Alamos National Laboratory
TA-00, SM1237, Rm104C
Los Alamos, New Mexico

Contact: Ms. Nita Patel

Workorder: 444566

Parmname	NOM	Sample	Qual	QC	Units	RPD%	REC%	Range	Anlst	Date	Time
Carbon Analysis											
Batch	1741854										
QC1203980265	444446006	DUP									
Total Organic Carbon Average		J	0.422	J	0.432	mg/L	2.34 ^	(+/-1.00)	TSM	02/28/18	22:29
QC1203980263	LCS										
Total Organic Carbon Average	10.0				10.5	mg/L		105 (80%-120%)		02/28/18	15:35
QC1203980262	MB										
Total Organic Carbon Average			U		ND	mg/L				02/28/18	15:25
QC1203980269	444446006	PS									
Total Organic Carbon Average	10.0	J	0.422		11.8	mg/L		114 (75%-125%)		02/28/18	23:09
Flow Injection Analysis											
Batch	1740518										
QC1203975197	444396002	DUP									
Cyanide, Total		U	ND	U	ND	ug/L	N/A		AXH3	02/28/18	07:57
QC1203975196	LCS										
Cyanide, Total	50.0				49.7	ug/L		99.4 (90%-110%)		02/28/18	07:41
QC1203975195	MB										
Cyanide, Total			U		ND	ug/L				02/28/18	07:40
QC1203975198	444396002	MS									
Cyanide, Total	100	U	ND		105	ug/L		105 (90%-110%)		02/28/18	07:58
Ion Chromatography											
Batch	1741949										
QC1203978238	444565001	DUP									
Bromide			0.210		0.209	mg/L	0.573 ^	(+/-0.200)	JXH5	02/23/18	19:21

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QC Summary

Workorder: 444566

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Parmname	NOM	Sample	Qual	QC	Units	RPD%	REC%	Range	Anlst	Date	Time
Ion Chromatography											
Batch	1741949										
Chloride		22.9		22.6	mg/L	1.2		(0%-20%)	JXH5	02/26/18	11:44
Fluoride		0.259		0.252	mg/L	2.74	^	(+/-0.100)		02/23/18	19:21
Sulfate		31.5		31.6	mg/L	0.233		(0%-20%)		02/26/18	11:44
QC1203978237	LCS										
Bromide	1.25			1.20	mg/L		95.8	(80%-120%)		02/23/18	18:19
Chloride	5.00			4.85	mg/L		97	(80%-120%)			
Fluoride	2.50			2.55	mg/L		102	(80%-120%)			
Sulfate	10.0			9.92	mg/L		99.2	(80%-120%)			
QC1203978236	MB										
Bromide			U	ND	mg/L					02/23/18	17:48
Chloride			U	ND	mg/L						
Fluoride			U	ND	mg/L						
Sulfate			U	ND	mg/L						
QC1203978239	444565001	PS									
Bromide	1.25	0.210		1.44	mg/L		98.3	(75%-125%)		02/23/18	19:52
Chloride	10.0	4.58		16.3	mg/L		117	(75%-125%)		02/26/18	12:15
Fluoride	2.50	0.259		2.72	mg/L		98.2	(75%-125%)		02/23/18	19:52
Sulfate	20.0	6.31		32.3	mg/L		130*	(75%-125%)		02/26/18	12:15

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QC Summary

Workorder: 444566

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Parmname	NOM	Sample	Qual	QC	Units	RPD%	REC%	Range	Anlst	Date	Time
Nutrient Analysis											
Batch	1741276										
QC1203976872	444566001	DUP									
Nitrogen, Ammonia		0.0759		0.0731	mg/L	3.76 ^		(+/-0.050)	KLP1	02/27/18	10:32
QC1203976870	LCS										
Nitrogen, Ammonia	1.00			0.938	mg/L		93.8	(90%-110%)		02/27/18	10:07
QC1203976869	MB										
Nitrogen, Ammonia			U	ND	mg/L					02/27/18	10:06
QC1203976874	444566001	MS									
Nitrogen, Ammonia	1.00	0.0759		1.03	mg/L		95.4	(90%-110%)		02/27/18	10:33
Batch	1742074										
QC1203978466	444396002	DUP									
Nitrogen, Total Kjeldahl		0.164	U	ND	mg/L	200* ^		(+/-0.100)	KLP1	02/28/18	14:14
QC1203978468	444396006	DUP									
Nitrogen, Total Kjeldahl		U	ND	U	ND	mg/L	N/A			02/28/18	14:53
QC1203978465	LCS										
Nitrogen, Total Kjeldahl	1.00			0.964	mg/L		96.4	(90%-110%)		02/28/18	14:13
QC1203978464	MB										
Nitrogen, Total Kjeldahl			U	ND	mg/L					02/28/18	14:12
QC1203978467	444396002	MS									
Nitrogen, Total Kjeldahl	1.00	0.164		0.665	mg/L		50.1 *	(90%-110%)		02/28/18	14:51
QC1203978469	444396006	MS									
Nitrogen, Total Kjeldahl	1.00	U	ND	0.751	mg/L		75.1 *	(90%-110%)		02/28/18	14:54
Batch	1742097										
QC1203978531	444396001	DUP									
Phosphorus, Total as P		0.0883		0.0898	mg/L	1.68 ^		(+/-0.050)	KLP1	03/05/18	12:06

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QC Summary

Workorder: 444566

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Parmname	NOM	Sample	Qual	QC	Units	RPD%	REC%	Range	Anlst	Date	Time
Nutrient Analysis											
Batch	1742097										
QC1203978530	LCS										
Phosphorus, Total as P	1.00			1.10	mg/L		110	(80%-124%)	KLP1	03/05/18	12:01
QC1203978529	MB										
Phosphorus, Total as P			U	ND	mg/L					03/05/18	12:00
QC1203978532	444396001	MS									
Phosphorus, Total as P	1.00	0.0883		1.09	mg/L		100	(63%-139%)		03/05/18	12:12
<hr/>											
Batch	1742109										
QC1203978569	444396001	DUP									
Nitrogen, Nitrate/Nitrite		0.783		0.781	mg/L	0.256		(0%-20%)	AXH3	02/26/18	09:01
QC1203978567	LCS										
Nitrogen, Nitrate/Nitrite	1.00			1.09	mg/L		109	(90%-110%)		02/26/18	08:48
QC1203978566	MB										
Nitrogen, Nitrate/Nitrite			U	ND	mg/L					02/26/18	08:47
QC1203978571	444396001	PS									
Nitrogen, Nitrate/Nitrite	1.00	0.783		1.74	mg/L		95.7	(90%-110%)		02/26/18	09:02
<hr/>											
Solids Analysis											
Batch	1742467										
QC1203979423	444446003	DUP									
Total Dissolved Solids		157		153	mg/L	2.76		(0%-5%)	KLP1	02/27/18	12:56
QC1203979422	LCS										
Total Dissolved Solids	300			293	mg/L		97.6	(95%-105%)		02/27/18	12:56
QC1203979421	MB										
Total Dissolved Solids			U	ND	mg/L					02/27/18	12:56

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QC Summary

Workorder: 444566

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Parmname	NOM	Sample	Qual	QC	Units	RPD%	REC%	Range	Anlst	Date	Time
Titration and Ion Analysis											
Batch	1741047										
QC1203976486	444026001	DUP									
Alkalinity, Total as CaCO3		58.0		57.4	mg/L	1.04		(0%-20%)	RXB5	02/24/18	14:29
Carbonate alkalinity (CaCO3)	U	ND	U	ND	mg/L	N/A					
QC1203976487	444396029	DUP									
Alkalinity, Total as CaCO3		65.2		64.4	mg/L	1.23		(0%-20%)		02/24/18	15:01
Carbonate alkalinity (CaCO3)	U	ND	U	ND	mg/L	N/A					
QC1203976485	LCS										
Alkalinity, Total as CaCO3	100			107	mg/L		107	(90%-110%)		02/24/18	14:24
QC1203976488	444026001	MS									
Alkalinity, Total as CaCO3	100	58.0		163	mg/L		105	(80%-120%)		02/24/18	14:30
QC1203976489	444396029	MS									
Alkalinity, Total as CaCO3	100	65.2		166	mg/L		101	(80%-120%)		02/24/18	15:03
Batch	1741049										
QC1203976495	444026001	DUP									
pH	H	7.93	H	7.91	SU	0.253		(0%-5%)	RXB5	02/24/18	14:27
QC1203976494	LCS										
pH	7.00			7.00	SU		100	(99%-101%)		02/24/18	14:25
Batch	1742552										
QC1203979623	444446001	DUP									
Conductivity		139		139	umhos/cm	0.144		(0%-10%)	HXC1	02/27/18	13:12
QC1203979622	LCS										
Conductivity	1410			1410	umhos/cm		99.6	(95%-105%)		02/27/18	13:11

Notes:

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QC Summary

Workorder: 444566

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Parmname	NOM	Sample	Qual	QC	Units	RPD%	REC%	Range	Anlst	Date	Time
<	Result is less than value reported										
>	Result is greater than value reported										
B	The target analyte was detected in the associated blank.										
E	General Chemistry--Concentration of the target analyte exceeds the instrument calibration range										
H	Analytical holding time was exceeded										
J	Value is estimated										
N/A	RPD or %Recovery limits do not apply.										
N1	See case narrative										
ND	Analyte concentration is not detected above the detection limit										
NJ	Consult Case Narrative, Data Summary package, or Project Manager concerning this qualifier										
Q	One or more quality control criteria have not been met. Refer to the applicable narrative or DER.										
R	Per section 9.3.4.1 of Method 1664 Revision B, due to matrix spike recovery issues, this result may not be reported or used for regulatory compliance purposes.										
R	Sample results are rejected										
U	Analyte was analyzed for, but not detected above the MDL, MDA, MDC or LOD.										
X	Consult Case Narrative, Data Summary package, or Project Manager concerning this qualifier										
Z	Paint Filter Test--Particulates passed through the filter, however no free liquids were observed.										
^	RPD of sample and duplicate evaluated using +/-RL. Concentrations are <5X the RL. Qualifier Not Applicable for Radiochemistry.										
d	5-day BOD--The 2:1 depletion requirement was not met for this sample										
e	5-day BOD--Test replicates show more than 30% difference between high and low values. The data is qualified per the method and can be used for reporting purposes										
h	Preparation or preservation holding time was exceeded										

N/A indicates that spike recovery limits do not apply when sample concentration exceeds spike conc. by a factor of 4 or more or %RPD not applicable.

^ The Relative Percent Difference (RPD) obtained from the sample duplicate (DUP) is evaluated against the acceptance criteria when the sample is greater than five times (5X) the contract required detection limit (RL). In cases where either the sample or duplicate value is less than 5X the RL, a control limit of +/- the RL is used to evaluate the DUP result.

* Indicates that a Quality Control parameter was not within specifications.

For PS, PSD, and SDILT results, the values listed are the measured amounts, not final concentrations.

Where the analytical method has been performed under NELAP certification, the analysis has met all of the requirements of the NELAC standard unless qualified on the QC Summary.

Radiological Analysis

Case Narrative

**Radiochemistry
Technical Case Narrative
ARS International, LLC (ARSL)
SDG #: 2018-1811
Work Order #: 444566**

Method/Analysis Information

Product: Alphaspec Am241 Liquid

Analytical Method: HASL-300:AM-241

Analytical Batch Number: 1741983

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566012	CAWA-18-56
1203978298	Method Blank (MB)
1203978300	Laboratory Control Sample (LCS)
1203978299	444566002(CAPA-18-2) Sample Duplicate (DUP)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-RAD-A-011 REV# 26.

Calibration Information:

Calibration Information

All initial and continuing calibration requirements have been met. Calibrations are performed monthly using mixed alpha standards comprised of the following: Gd-148, Np-237, and Cm-244. The initial Calibration was performed in February 2018.

Standards Information

Standard solutions for these analysis are NIST traceable or verified with a NIST traceable standard and used before the expiration dates.

Sample Geometry

All counting sources were prepared in the same geometry as the calibration standards.

Quality Control (QC) Information:

Blank Information

Aliquots for samples 1203978298 (MB) and 1203978300 (LCS) were changed to 1.0, and the MDCs (and Lc if requested) for all samples are calculated using a blank population per client request.

Method Blank Criteria

The Method Blank (MB) met acceptance criteria.

CSU

The blank result is less than 1.65 times the CSU.

Blank Decision Level

The blank result is less than the decision level.

Tracer/Carrier Yield

All yields met the required acceptance limits.

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries met the acceptance limits.

Designated QC

The following sample was used for QC: 444566002 (CAPA-18-2). The QC was from ARSL work order 444566.

Duplication Criteria between QC Sample and Duplicate Sample

The QC Sample and Duplicate Sample (DUP) met the duplication acceptance criteria.

RDL Met

The method RDL has been met.

Technical Information:**Holding Time**

All sample procedures for this sample set were performed within the required holding time.

Negative > 3 sigma TPU

Samples results are not more negative than three sigma TPU.

Sample Re-prep/Re-analysis

None of the samples in this sample set required prep or reanalysis.

Recounts

None of the samples in this sample set were recounted.

Miscellaneous Information:**Manual Integration**

No manual integrations were performed on data in this batch.

Sample-Specific MDA/MDC

The MDA/MDC reported on the certificate of analysis is a sample-specific MDA/MDC.

Additional Comments

Additional comments were not required for this sample set.

Qualifier Information

Manual qualifiers were not required.

Method/Analysis Information

Product: ISOPU
Analytical Method: HASL-300:ISOPU
Analytical Batch Number: 1741984

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566012	CAWA-18-56
1203978301	Method Blank (MB)
1203978303	Laboratory Control Sample (LCS)
1203978302	444566002(CAPA-18-2) Sample Duplicate (DUP)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-RAD-A-011 REV# 26.

Calibration Information:

Calibration Information

All initial and continuing calibration requirements have been met. Calibrations are performed monthly using mixed alpha standards comprised of the following: Gd-148, Np-237, and Cm-244. The initial Calibration was performed in February 2018.

Standards Information

Standard solutions for these analysis are NIST traceable or verified with a NIST traceable standard and used before the expiration dates.

Sample Geometry

All counting sources were prepared in the same geometry as the calibration standards.

Quality Control (QC) Information:

Blank Information

Aliquots for samples 1203978301 (MB) and 1203978303 (LCS) were changed to 1.0, and the MDCs (and Lc if requested) for all samples are calculated using a blank population per client request.

Method Blank Criteria

The Method Blank (MB) met acceptance criteria.

CSU

The blank result is less than 1.65 times the CSU.

Blank Decision Level

The blank result is less than the decision level.

Tracer/Carrier Yield

All yields met the required acceptance limits.

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries met the acceptance limits.

Designated QC

The following sample was used for QC: 444566002 (CAPA-18-2). The QC was from ARSL work order 444566.

Duplication Criteria between QC Sample and Duplicate Sample

The QC Sample and Duplicate Sample (DUP) met the duplication acceptance criteria.

RDL Met

Sample (See Below) did not meet the detection limit due to the high standard deviation. When a blank population is performed the MDC is greater than the RDL due to the high standard deviation. The sample was counted the maximum count time of in order to achieve the lowest possible MDAs.

Sample	Analyte	Value
444566002 (CAPA-18-2)	Plutonium-238	Result 0.00332 < MDA 0.0577 > RDL 0.05 pCi/L

Technical Information:**Holding Time**

All sample procedures for this sample set were performed within the required holding time.

Negative > 3 sigma TPU

Samples results are not more negative than three sigma TPU.

Sample Re-prep/Re-analysis

None of the samples in this sample set required reprep or reanalysis.

Recounts

None of the samples in this sample set were recounted.

Miscellaneous Information:**Manual Integration**

No manual integrations were performed on data in this batch.

Sample-Specific MDA/MDC

The MDA/MDC reported on the certificate of analysis is a sample-specific MDA/MDC.

Additional Comments

Additional comments were not required for this sample set.

Qualifier Information

Manual qualifiers were not required.

Method/Analysis Information

Product: IsoU

Analytical Method: HASL-300:ISOU

Analytical Batch Number: 1741985

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566012	CAWA-18-56
1203978305	Method Blank (MB)
1203978307	Laboratory Control Sample (LCS)
1203978306	444566002(CAPA-18-2) Sample Duplicate (DUP)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-RAD-A-011 REV# 26.

Calibration Information:

Calibration Information

All initial and continuing calibration requirements have been met. Calibrations are performed monthly using mixed alpha standards comprised of the following: Gd-148, Np-237, and Cm-244. The initial Calibration was performed in February 2018.

Standards Information

Standard solutions for these analysis are NIST traceable or verified with a NIST traceable standard and used before the expiration dates.

Sample Geometry

All counting sources were prepared in the same geometry as the calibration standards.

Quality Control (QC) Information:

Blank Information

Aliquots for samples 1203978305 (MB) and 1203978307 (LCS) were changed to 1.0, and the MDCs (and Lc if requested) for all samples are calculated using a blank population per client request.

Method Blank Criteria

The Method Blank (MB) met acceptance criteria.

CSU

The blank (See Below) result is greater than 1.65 times the CSU but less than the MDC.

Sample	Analyte	Value
1203978305 (MB)	Uranium-235/236	Blank result > 1.65 CSU

Blank Decision Level

The blank result is less than the decision level.

Tracer/Carrier Yield

All yields met the required acceptance limits.

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries met the acceptance limits.

Designated QC

The following sample was used for QC: 444566002 (CAPA-18-2). The QC was from ARSL work order 444566.

Duplication Criteria between QC Sample and Duplicate Sample

The QC Sample and Duplicate Sample (DUP) met the duplication acceptance criteria.

RDL Met

The method RDL has been met.

Technical Information:**Holding Time**

All sample procedures for this sample set were performed within the required holding time.

Negative > 3 sigma TPU

Samples results are not more negative than three sigma TPU.

Sample Re-prep/Re-analysis

None of the samples in this sample set required prep or reanalysis.

Recounts

Samples 1203978306 (CAPA-18-2DUP) and 444566012 (CAWA-18-56) were recounted due to a peak shift.

The recounts are reported.

Miscellaneous Information:**Manual Integration**

No manual integrations were performed on data in this batch.

Sample-Specific MDA/MDC

The MDA/MDC reported on the certificate of analysis is a sample-specific MDA/MDC.

Additional Comments

Additional comments were not required for this sample set.

Qualifier Information

Manual qualifiers were not required.

Method/Analysis Information

Product: **Gammaspec**

Analytical Method: EPA:901.1

Analytical Batch Number: 1741844

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566012	CAWA-18-56
1203978021	Method Blank (MB)
1203978023	Laboratory Control Sample (LCS)
1203978022	444566002(CAPA-18-2) Sample Duplicate (DUP)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-RAD-A-013 REV# 27.

Calibration Information:

Calibration Information

All initial and continuing calibration requirements have been met. The initial Calibrations were performed in August 2017, January 2018, June 2017, May 2017 and November 2017.

Standards Information

Standard solutions for these analysis are NIST traceable or verified with a NIST traceable standard and used before the expiration dates.

Sample Geometry

All counting sources were prepared in the same geometry as the calibration standards.

Quality Control (QC) Information:

Blank Information

The blank volume is representative of the sample volume in this batch.

Method Blank Criteria

The Method Blank (MB) met acceptance criteria.

CSU

The blank (See Below) result is greater than 1.65 times the CSU but less than the MDC.

Sample	Analyte	Value
1203978021 (MB)	Cobalt-60	Blank result > 1.65 CSU

Blank Decision Level

The blank (See Below) result is greater than the decision level but less than the MDC.

Sample	Analyte	Value
1203978021 (MB)	Cobalt-60	Blank result > DL

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries met the acceptance limits.

Designated QC

The following sample was used for QC: 444566002 (CAPA-18-2). The QC was from ARSL work order 444566.

Duplication Criteria between QC Sample and Duplicate Sample

The QC Sample and Duplicate Sample (DUP) met the duplication acceptance criteria.

RDL Met

The method RDL has been met.

Technical Information:

Holding Time

All sample procedures for this sample set were performed within the required holding time.

Sample Re-prep/Re-analysis

None of the samples in this sample set required reprep or reanalysis.

Recounts

None of the samples in this sample set were recounted.

Additional Identified Radionuclides

No additional radionuclides were added.

Miscellaneous Information:

Sample-Specific MDA/MDC

The MDA/MDC reported on the certificate of analysis is a sample-specific MDA/MDC.

Additional Comments

Additional comments were not required for this sample set.

Qualifier Information

Manual qualifiers were not required.

Method/Analysis Information

Product:	WSP-GrossA/B
Analytical Method:	EPA 900.0/SW846 9310
Analytical Batch Number:	1742280

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566012	CAWA-18-56
1203978996	Method Blank (MB)
1203979000	Laboratory Control Sample (LCS)
1203978997	444396013(CAWA-18-40) Sample Duplicate (DUP)
1203978998	444396013(CAWA-18-40) Matrix Spike (MS)
1203978999	444396013(CAWA-18-40) Matrix Spike Duplicate (MSD)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-RAD-A-001 REV# 19.

Calibration Information:

Calibration Information

All initial and continuing calibration requirements have been met. The initial Calibration was performed in December 2017.

Standards Information

Standard solutions for these analysis are NIST traceable or verified with a NIST traceable standard and used before the expiration dates.

Sample Geometry

All counting sources were prepared in the same geometry as the calibration standards.

Quality Control (QC) Information:

Blank Information

Aliquots for samples 1203978996 (MB) and 1203979000 (LCS) were changed to 1.0 per client request.

Method Blank Criteria

The Method Blank (MB) met acceptance criteria.

CSU

The blank result is less than 1.65 times the CSU.

Blank Decision Level

The blank result is less than the decision level.

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries met the acceptance limits.

Designated QC

The following sample was used for QC: 444396013 (CAWA-18-40). The QC was from ARSL work order 444396.

Matrix Spike (MS) Recovery

The MS spike recoveries met acceptance limits.

Duplication Criteria between MS and MSD

The Matrix Spike (MS) and Matrix Spike Duplicate (MSD) met the duplication acceptance criteria.

Duplication Criteria between QC Sample and Duplicate Sample

The QC Sample and Duplicate Sample (DUP) met the duplication acceptance criteria.

RDL Met

The method RDL has been met.

Technical Information:**Holding Time**

All sample procedures for this sample set were performed within the required holding time.

Negative > 3 sigma TPU

Samples results are not more negative than three sigma TPU.

Sample Re-prep/Re-analysis

None of the samples in this sample set required reprep or reanalysis.

Gross Alpha/Beta Preparation Information

High hygroscopic salt content in evaporated samples can cause the sample mass to fluctuate due to moisture absorption. To minimize this interference, the salts are converted to oxides by heating the sample under a flame until a dull red color is obtained. The conversion to oxides stabilizes the sample weight and ensures that proper alpha/beta efficiencies are assigned for each sample. Volatile radioisotopes of carbon, hydrogen, technetium, polonium and cesium may be lost during sample heating, especially to a dull red heat. For this sample set, the prepared planchet was counted for beta activity before being flamed. After flaming, the planchet was counted for alpha activity.

Recounts

None of the samples in this sample set were recounted.

Miscellaneous Information:**Sample-Specific MDA/MDC**

The MDA/MDC reported on the certificate of analysis is a sample-specific MDA/MDC.

Additional Comments

The matrix spike and matrix spike duplicate, 1203978998 (CAWA-18-40MS) and 1203978999 (CAWA-18-40MSD), aliquots were reduced to conserve sample volume.

Qualifier Information

Manual qualifiers were not required.

Method/Analysis Information

Product: GFPC, Sr90, liquid

Analytical Method: EPA:905.0

Analytical Batch Number: 1742296

Sample ID	Client ID
444566002	CAPA-18-2
444566006	CAPA-18-12
444566012	CAWA-18-56
1203979041	Method Blank (MB)
1203979044	Laboratory Control Sample (LCS)
1203979042	444617006(CAWA-18-67) Sample Duplicate (DUP)
1203979043	444617006(CAWA-18-67) Matrix Spike (MS)

The samples in this SDG were analyzed on an "as received" basis.

SOP Reference

Procedure for preparation, analysis and reporting of analytical data are controlled by GEL Laboratories LLC as Standard Operating Procedure (SOP). The data discussed in this narrative has been analyzed in accordance with GL-RAD-A-004 REV# 19.

Calibration Information:

Calibration Information

All initial and continuing calibration requirements have been met. The initial Calibration was performed in April 2016.

Standards Information

Standard solutions for these analysis are NIST traceable or verified with a NIST traceable standard and used before the expiration dates.

Sample Geometry

All counting sources were prepared in the same geometry as the calibration standards.

Quality Control (QC) Information:

Blank Information

Aliquots for samples 1203979041 (MB) and 1203979044 (LCS) were changed to 1.0 per client request.

Method Blank Criteria

The Method Blank (MB) met acceptance criteria.

CSU

The blank result is less than 1.65 times the CSU.

Blank Decision Level

The blank result is less than the decision level.

Tracer/Carrier Yield

All yields met the required acceptance limits.

Laboratory Control Sample (LCS) Recovery

The LCS spike recoveries met the acceptance limits.

Designated QC

The following sample was used for QC: 444617006 (CAWA-18-67). The QC was from ARSL work order 444617.

Matrix Spike (MS) Recovery

The MS spike recoveries met acceptance limits.

Duplication Criteria between QC Sample and Duplicate Sample

The QC Sample and Duplicate Sample (DUP) met the duplication acceptance criteria.

RDL Met

The method RDL has been met.

Technical Information:**Holding Time**

All sample procedures for this sample set were performed within the required holding time.

Negative > 3 sigma TPU

Samples results are not more negative than three sigma TPU.

Sample Re-prep/Re-analysis

None of the samples in this sample set required reprep or reanalysis.

Recounts

Sample 1203979041 (MB) was recounted due to a suspected blank false positive. The recount is reported.
Sample 444566002 (CAPA-18-2) was recounted due to a suspected false positive. The recount is reported.

Miscellaneous Information:**Sample-Specific MDA/MDC**

The MDA/MDC reported on the certificate of analysis is a sample-specific MDA/MDC.

Additional Comments

The matrix spike, 1203979043 (CAWA-18-67MS), aliquot was reduced to conserve sample volume.

Qualifier Information

Manual qualifiers were not required.

Certification Statement

Where the analytical method has been performed under NELAP certification, the analysis has met all of the requirements of the NELAC standard unless otherwise noted in the analytical case narrative.

GEL LABORATORIES LLC

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Qualifier Definition Report for

ARSL004 ARS International, LLC (ARS-LANS-MTOA6-25093-GEL)

Client SDG: 2018-1811 GEL Work Order: 444566


The Qualifiers in this report are defined as follows:

- * A quality control analyte recovery is outside of specified acceptance criteria
- ** Analyte is a Tracer compound
- U Analyte was analyzed for, but not detected above the MDL, MDA, MDC or LOD.

Review/Validation

GEL requires all analytical data to be verified by a qualified data reviewer. In addition, all CLP-like deliverables receive a third level review of the fractional data package.

The following data validator verified the information presented in this data report:

Signature: 

Name: Theresa Austin

Date: 12 MAR 2018

Title: Group Leader

Sample Data Summary

GEL LABORATORIES LLC

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Certificate of Analysis

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545

Contact: Ms. Nita Patel

Project: LANL- WQH Water Samples

Report Date: March 12, 2018

Client Sample ID: CAPA-18-2
Sample ID: 444566002
Matrix: W
Collect Date: 21-FEB-18
Receive Date: 23-FEB-18
Collector: Client

Project: ESHL00114
Client ID: ARSL004

Parameter	Qualifier	Result	Uncertainty	MDC	Lc	TPU	RL	Units	PF	DF	Analyst	Date	Time	Batch	Mtd.
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Rad Alpha Spec Analysis

Alphaspec Am241 Liquid "As Received"

Americium-241	U	0.0071	+/-0.00561	0.030	0.0126	+/-0.00562	0.050	pCi/L			MXS2	02/27/18	1357	1741983	1
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ISOPU "As Received"

Plutonium-238	U	0.00332	+/-0.00742	0.0577	0.0243	+/-0.00742	0.050	pCi/L			MXS2	02/27/18	1357	1741984	2
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Plutonium-239/240	U	0.0166	+/-0.011	0.0464	0.0187	+/-0.011	0.050	pCi/L							
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IsoU "As Received"

Uranium-234		0.291	+/-0.0286	0.112	0.0528	+/-0.0321	1.00	pCi/L			MXS2	02/27/18	1332	1741985	3
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Uranium-235/236	U	0.0209	+/-0.0123	0.0625	0.0272	+/-0.0123	1.00	pCi/L							
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Uranium-238		0.191	+/-0.0235	0.0567	0.0251	+/-0.0253	0.500	pCi/L							
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Rad Gamma Spec Analysis

Gammasespec "As Received"

Cesium-137	U	-0.858	+/-1.14	3.61	1.58	+/-1.16	8.00	pCi/L			BSW1	02/23/18	1126	1741844	4
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Cobalt-60	U	0.0705	+/-1.11	4.29	1.80	+/-1.11	8.00	pCi/L							
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Neptunium-237	U	-0.856	+/-1.85	6.58	3.02	+/-1.86		pCi/L							
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Potassium-40	U	-19.9	+/-16.7	58.0	25.5	+/-17.3		pCi/L							
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Sodium-22	U	-2.1	+/-1.13	3.40	1.36	+/-1.23		pCi/L							
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Rad Gas Flow Proportional Counting

GFPC, Sr90, liquid "As Received"

Strontium-90	U	-0.168	+/-0.131	0.482	0.226	+/-0.131	0.500	pCi/L			KSD1	03/01/18	0836	1742296	5
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WSP-GrossA/B "As Received"

Beta		3.26	+/-0.801	2.33	1.06	+/-0.846	3.00	pCi/L			BXG2	03/08/18	0830	1742280	6
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Alpha	U	0.947	+/-0.785	2.78	0.981	+/-0.790	3.00	pCi/L			BXG2	03/08/18	1338	1742280	7
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The following Analytical Methods were performed

Method	Description
1	HASL-300:AM-241
2	HASL-300:ISOPU
3	HASL-300:ISOU
4	EPA:901.1
5	EPA:905.0
6	EPA 900.0/SW846 9310
7	EPA 900.0/SW846 9310

Surrogate/Tracer Recovery	Test	Batch ID	Recovery%	Acceptable Limits
Americium-243 Tracer	Alphaspec Am241 Liquid "As Received"	1741983	93.5	(50%-105%)
Plutonium-242 Tracer	ISOPU "As Received"	1741984	60.3	(50%-105%)
Uranium-232 Tracer	IsoU "As Received"	1741985	69	(50%-105%)

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Certificate of Analysis

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545

Contact: Ms. Nita Patel

Project: LANL- WQH Water Samples

Client Sample ID: CAPA-18-2

Sample ID: 444566002

Project: ESHL00114

Client ID: ARSL004

Report Date: March 12, 2018

Parameter	Qualifier	Result	Uncertainty	MDC	Lc	TPU	RL	Units	PF	DF	Analyst	Date	Time	Batch	Mtd.
Surrogate/Tracer Recovery	Test								Batch ID	Recovery%	Acceptable Limits				
Strontium Carrier		GFPC, Sr90, liquid "As Received"							1742296	97.7	(50%-105%)				

Notes:

The MDC is a sample specific MDC.

TPU and Counting Uncertainty are calculated at the 68% confidence level (1-sigma).

Column headers are defined as follows:

DF: Dilution Factor

DL: Detection Limit

Lc/LC: Critical Level

MDA: Minimum Detectable Activity

MDC: Minimum Detectable Concentration

Mtd.: Method

PF: Prep Factor

RL: Reporting Limit

TPU: Total Propagated Uncertainty

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Certificate of Analysis

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545

Contact: Ms. Nita Patel

Project: LANL- WQH Water Samples

Client Sample ID: CAPA-18-12

Sample ID: 444566006

Matrix: W

Collect Date: 21-FEB-18

Receive Date: 23-FEB-18

Collector: Client

Report Date: March 12, 2018

Project: ESHL00114

Client ID: ARSL004

Parameter	Qualifier	Result	Uncertainty	MDC	Lc	TPU	RL	Units	PF	DF	Analyst	Date	Time	Batch	Mtd.
Rad Alpha Spec Analysis															
<i>Alphaspec Am241 Liquid "As Received"</i>															
Americium-241	U	0.00371	+/-0.00586	0.0313	0.0132	+/-0.00586	0.050	pCi/L			MXS2	02/27/18	1357	1741983	1
<i>ISOPU "As Received"</i>															
Plutonium-238	U	0.00425	+/-0.00672	0.0369	0.0156	+/-0.00672	0.050	pCi/L			MXS2	02/27/18	1357	1741984	2
Plutonium-239/240	U	0.00213	+/-0.00766	0.0297	0.012	+/-0.00766	0.050	pCi/L							
<i>IsoU "As Received"</i>															
Uranium-234		0.277	+/-0.0288	0.115	0.0542	+/-0.032	1.00	pCi/L			MXS2	02/27/18	1332	1741985	3
Uranium-235/236	U	0.0153	+/-0.00811	0.0641	0.0279	+/-0.00814	1.00	pCi/L							
Uranium-238		0.193	+/-0.0235	0.0582	0.0257	+/-0.0253	0.500	pCi/L							
Rad Gamma Spec Analysis															
<i>Gammasespec "As Received"</i>															
Cesium-137	U	-0.267	+/-0.935	3.35	1.49	+/-0.937	8.00	pCi/L			BSW1	02/23/18	1140	1741844	4
Cobalt-60	U	0.428	+/-0.794	3.36	1.40	+/-0.800	8.00	pCi/L							
Neptunium-237	U	-0.991	+/-1.82	6.54	3.02	+/-1.83		pCi/L							
Potassium-40	U	-19.6	+/-13.5	47.8	21.2	+/-14.3		pCi/L							
Sodium-22	U	-0.544	+/-0.680	2.54	1.00	+/-0.692		pCi/L							
Rad Gas Flow Proportional Counting															
<i>GFPC, Sr90, liquid "As Received"</i>															
Strontium-90	U	0.016	+/-0.0812	0.278	0.133	+/-0.0812	0.500	pCi/L			KSD1	02/28/18	1617	1742296	5
<i>WSP-GrossA/B "As Received"</i>															
Beta		4.76	+/-0.895	2.46	1.12	+/-0.978	3.00	pCi/L			BXG2	03/08/18	0830	1742280	6
Alpha	U	1.07	+/-0.781	2.66	0.897	+/-0.787	3.00	pCi/L			BXG2	03/08/18	1338	1742280	7

The following Analytical Methods were performed

Method	Description
1	HASL-300:AM-241
2	HASL-300:ISOPU
3	HASL-300:ISOU
4	EPA:901.1
5	EPA:905.0
6	EPA 900.0/SW846 9310
7	EPA 900.0/SW846 9310

Surrogate/Tracer	Recovery	Test	Batch ID	Recovery%	Acceptable Limits
Americium-243 Tracer		Alphaspec Am241 Liquid "As Received"	1741983	102	(50%-105%)
Plutonium-242 Tracer		ISOPU "As Received"	1741984	83.9	(50%-105%)
Uranium-232 Tracer		IsoU "As Received"	1741985	76.4	(50%-105%)
Strontium Carrier		GFPC, Sr90, liquid "As Received"	1742296	81.4	(50%-105%)

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Certificate of Analysis

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545

Contact: Ms. Nita Patel

Project: LANL- WQH Water Samples

Client Sample ID: CAPA-18-12

Sample ID: 444566006

Project: ESHL00114

Client ID: ARSL004

Report Date: March 12, 2018

Parameter	Qualifier	Result	Uncertainty	MDC	Lc	TPU	RL	Units	PF	DF	Analyst	Date	Time	Batch	Mtd.
Surrogate/Tracer Recovery	Test											Batch ID	Recovery%	Acceptable Limits	

Notes:

The MDC is a sample specific MDC.

TPU and Counting Uncertainty are calculated at the 68% confidence level (1-sigma).

Column headers are defined as follows:

DF: Dilution Factor

DL: Detection Limit

Lc/LC: Critical Level

MDA: Minimum Detectable Activity

MDC: Minimum Detectable Concentration

Mtd.: Method

PF: Prep Factor

RL: Reporting Limit

TPU: Total Propagated Uncertainty

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Certificate of Analysis

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545

Contact: Ms. Nita Patel

Project: LANL- WQH Water Samples

Client Sample ID: CAWA-18-56

Sample ID: 444566012

Matrix: W

Collect Date: 21-FEB-18

Receive Date: 23-FEB-18

Collector: Client

Report Date: March 12, 2018

Project: ESHL00114

Client ID: ARSL004

Parameter	Qualifier	Result	Uncertainty	MDC	Lc	TPU	RL	Units	PF	DF	Analyst	Date	Time	Batch	Mtd.
Rad Alpha Spec Analysis															
<i>Alphaspec Am241 Liquid "As Received"</i>															
Americium-241	U	0.00213	+/-0.00369	0.036	0.0151	+/-0.00369	0.050	pCi/L			MXS2	02/27/18	1357	1741983	1
<i>ISOPU "As Received"</i>															
Plutonium-238	U	-0.0065	+/-0.00782	0.0377	0.0159	+/-0.00782	0.050	pCi/L			MXS2	02/27/18	1357	1741984	2
Plutonium-239/240	U	0.0065	+/-0.0108	0.0303	0.0122	+/-0.0108	0.050	pCi/L							
<i>IsoU "As Received"</i>															
Uranium-234		0.239	+/-0.0314	0.153	0.0721	+/-0.0339	1.00	pCi/L			MXS2	03/05/18	1516	1741985	3
Uranium-235/236	U	0.0244	+/-0.0129	0.0853	0.0371	+/-0.0129	1.00	pCi/L							
Uranium-238		0.0989	+/-0.0198	0.0774	0.0342	+/-0.0204	0.500	pCi/L							
Rad Gamma Spec Analysis															
<i>Gammasespec "As Received"</i>															
Cesium-137	U	0.577	+/-0.918	3.51	1.57	+/-0.928	8.00	pCi/L			BSW1	02/23/18	1540	1741844	4
Cobalt-60	U	-1.13	+/-0.961	3.37	1.41	+/-0.997	8.00	pCi/L							
Neptunium-237	U	0.767	+/-2.40	7.56	3.53	+/-2.40		pCi/L							
Potassium-40	U	27.8	+/-20.4	38.3	16.4	+/-20.5		pCi/L							
Sodium-22	U	0.267	+/-0.951	3.82	1.64	+/-0.953		pCi/L							
Rad Gas Flow Proportional Counting															
<i>GFPC, Sr90, liquid "As Received"</i>															
Strontium-90	U	-0.0293	+/-0.0694	0.241	0.115	+/-0.0694	0.500	pCi/L			KSD1	02/28/18	1617	1742296	5
<i>WSP-GrossA/B "As Received"</i>															
Beta		3.38	+/-0.808	2.35	1.07	+/-0.855	3.00	pCi/L			BXG2	03/08/18	0830	1742280	6
Alpha	U	0.700	+/-0.765	2.76	1.12	+/-0.768	3.00	pCi/L			BXG2	03/08/18	1337	1742280	7

The following Analytical Methods were performed

Method	Description
1	HASL-300:AM-241
2	HASL-300:ISOPU
3	HASL-300:ISOU
4	EPA:901.1
5	EPA:905.0
6	EPA 900.0/SW846 9310
7	EPA 900.0/SW846 9310

Surrogate/Tracer	Recovery	Test	Batch ID	Recovery%	Acceptable Limits
Americium-243 Tracer		Alphaspec Am241 Liquid "As Received"	1741983	96	(50%-105%)
Plutonium-242 Tracer		ISOPU "As Received"	1741984	80.4	(50%-105%)
Uranium-232 Tracer		IsoU "As Received"	1741985	77.2	(50%-105%)
Strontium Carrier		GFPC, Sr90, liquid "As Received"	1742296	95.3	(50%-105%)

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Certificate of Analysis

Company : Los Alamos National Laboratory
Address : TA-00, SM1237, Rm104C

Los Alamos, New Mexico 87545

Contact: Ms. Nita Patel

Project: LANL- WQH Water Samples

Client Sample ID: CAWA-18-56

Sample ID: 444566012

Project: ESHL00114

Client ID: ARSL004

Report Date: March 12, 2018

Parameter	Qualifier	Result	Uncertainty	MDC	Lc	TPU	RL	Units	PF	DF	Analyst	Date	Time	Batch	Mtd.
Surrogate/Tracer Recovery	Test											Batch ID	Recovery%	Acceptable Limits	

Notes:

The MDC is a sample specific MDC.

TPU and Counting Uncertainty are calculated at the 68% confidence level (1-sigma).

Column headers are defined as follows:

DF: Dilution Factor

DL: Detection Limit

Lc/LC: Critical Level

MDA: Minimum Detectable Activity

MDC: Minimum Detectable Concentration

Mtd.: Method

PF: Prep Factor

RL: Reporting Limit

TPU: Total Propagated Uncertainty

Quality Control Summary

GEL LABORATORIES LLC

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QC Summary

Report Date: March 12, 2018

Page 1 of 6

Client : Los Alamos National Laboratory
TA-00, SM1237, Rm104C

Los Alamos, New Mexico

Contact: Ms. Nita Patel

Workorder: 444566

Parmname	NOM	Sample	Qual	QC	Units	RER	REC%	Range	Anlst	Date	Time
Rad Alpha Spec											
Batch	1741983										
QC1203978299	444566002	DUP									
Americium-241	U	0.0071	U	0.0035	pCi/L	0.148		(0-1)	MXS2	02/27/18	13:57
	Uncert:	+/-0.00561		+/-0.00655							
	TPU:	+/-0.00562		+/-0.00655							
**Americium-243 Tracer	2.60	2.45		2.57	pCi/L		99.1	(50%-105%)			
	Uncert:	+/-0.068		+/-0.0671							
	TPU:	+/-0.130		+/-0.129							
QC1203978300	LCS										
Americium-241	1.97			2.02	pCi/L		102	(80%-120%)	MXS2	02/27/18	13:56
	Uncert:			+/-0.0581							
	TPU:			+/-0.106							
**Americium-243 Tracer	2.10			2.19	pCi/L		104	(50%-105%)			
	Uncert:			+/-0.0588							
	TPU:			+/-0.109							
QC1203978298	MB										
Americium-241			U	0.0026	pCi/L				MXS2	02/27/18	13:57
	Uncert:			+/-0.00367							
	TPU:			+/-0.00367							
**Americium-243 Tracer	2.10			2.19	pCi/L		105	(50%-105%)			
	Uncert:			+/-0.052							
	TPU:			+/-0.102							
Batch	1741984										
QC1203978302	444566002	DUP									
Plutonium-238	U	0.00332	U	0.00431	pCi/L	0.0347		(0-1)	MXS2	02/27/18	13:57
	Uncert:	+/-0.00742		+/-0.00681							
	TPU:	+/-0.00742		+/-0.00681							
Plutonium-239/240	U	0.0166	U	0.00215	pCi/L	0.384		(0-1)			
	Uncert:	+/-0.011		+/-0.00776							
	TPU:	+/-0.011		+/-0.00776							
**Plutonium-242 Tracer	2.45	1.49		2.07	pCi/L		84.4	(50%-105%)			
	Uncert:	+/-0.0914		+/-0.073							
	TPU:	+/-0.147		+/-0.124							
QC1203978303	LCS										
Plutonium-238			U	0.00879	pCi/L			(80%-120%)	MXS2	02/27/18	13:56
	Uncert:			+/-0.00695							
	TPU:			+/-0.00696							
Plutonium-239/240	1.98			2.30	pCi/L		116	(80%-120%)			
	Uncert:			+/-0.0711							
	TPU:			+/-0.123							
**Plutonium-242 Tracer	1.98			1.60	pCi/L		80.6	(50%-105%)			
	Uncert:			+/-0.0664							
	TPU:			+/-0.109							

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QC Summary

Workorder: 444566

Page 2 of 6

Parmname	NOM	Sample	Qual	QC	Units	RER	REC%	Range	Anlst	Date	Time
Rad Alpha Spec											
Batch	1741984										
QC1203978301	MB										
Plutonium-238			U	0.00347	pCi/L				MXS2	02/27/18	13:57
				Uncert: +/-0.00649							
				TPU: +/-0.00649							
Plutonium-239/240			U	0.00347	pCi/L						
				Uncert: +/-0.00649							
				TPU: +/-0.00649							
**Plutonium-242 Tracer	1.98			1.45	pCi/L		73.5	(50%-105%)			
				Uncert: +/-0.0591							
				TPU: +/-0.100							
Batch	1741985										
QC1203978306	444566002	DUP									
Uranium-234		0.291		0.287	pCi/L	0.0342		(0-1)	MXS2	03/05/18	15:16
		Uncert: +/-0.0286		+/-0.0304							
		TPU: +/-0.0321		+/-0.0338							
Uranium-235/236		U 0.0209	U	0.0248	pCi/L	0.0802		(0-1)			
		Uncert: +/-0.0123		+/-0.0117							
		TPU: +/-0.0123		+/-0.0118							
Uranium-238		0.191		0.218	pCi/L	0.249		(0-1)			
		Uncert: +/-0.0235		+/-0.0262							
		TPU: +/-0.0253		+/-0.0284							
**Uranium-232 Tracer	2.59	1.80		1.98	pCi/L		76.4	(50%-105%)			
		Uncert: +/-0.082		+/-0.0867							
		TPU: +/-0.151		+/-0.156							
QC1203978307	LCS										
Uranium-234				2.80	pCi/L				MXS2	02/27/18	13:32
		Uncert: +/-0.0812		+/-0.0812							
		TPU: +/-0.163		+/-0.163							
Uranium-235/236				0.158	pCi/L						
		Uncert: +/-0.0225		+/-0.0225							
		TPU: +/-0.0238		+/-0.0238							
Uranium-238	2.70			2.95	pCi/L		109	(80%-120%)			
		Uncert: +/-0.0831		+/-0.0831							
		TPU: +/-0.170		+/-0.170							
**Uranium-232 Tracer	2.09			1.61	pCi/L		77.3	(50%-105%)			
		Uncert: +/-0.0703		+/-0.0703							
		TPU: +/-0.126		+/-0.126							
QC1203978305	MB										
Uranium-234			U	0.0106	pCi/L				MXS2	02/27/18	13:32
		Uncert: +/-0.00967		+/-0.00967							
		TPU: +/-0.00971		+/-0.00971							
Uranium-235/236			U	0.0187	pCi/L						
		Uncert: +/-0.00886		+/-0.00886							
		TPU: +/-0.00891		+/-0.00891							
Uranium-238			U	0.00432	pCi/L						
		Uncert: +/-0.00612		+/-0.00612							
		TPU: +/-0.00612		+/-0.00612							

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Parmname	NOM	Sample	Qual	QC	Units	RER	REC%	Range	Anlst	Date	Time
Rad Alpha Spec											
Batch	1741985										
*Uranium-232 Tracer	2.09			1.68	pCi/L		80.7	(50%-105%)			
	Uncert:			+/-0.0679							
	TPU:			+/-0.124							
Rad Gamma Spec											
Batch	1741844										
QC1203978022	444566002	DUP									
Cesium-137	U	-0.858	U	1.94	pCi/L	0.548		(0-1)	BSW1	02/23/18	15:41
	Uncert:	+/-1.14		+/-1.31							
	TPU:	+/-1.16		+/-1.39							
Cobalt-60	U	0.0705	U	-1.5	pCi/L	0.329		(0-1)			
	Uncert:	+/-1.11		+/-1.23							
	TPU:	+/-1.11		+/-1.28							
Neptunium-237	U	-0.856	U	-3.28	pCi/L	0.248		(0-1)			
	Uncert:	+/-1.85		+/-2.92							
	TPU:	+/-1.86		+/-3.02							
Potassium-40	U	-19.9	U	-46	pCi/L	0.311		(0-1)			
	Uncert:	+/-16.7		+/-22.0							
	TPU:	+/-17.3		+/-24.5							
Sodium-22	U	-2.1	U	0.255	pCi/L	0.4		(0-1)			
	Uncert:	+/-1.13		+/-1.71							
	TPU:	+/-1.23		+/-1.71							
QC1203978023	LCS										
Americium-241	34300			35800	pCi/L		104	(80%-120%)	BSW1	02/23/18	14:56
	Uncert:			+/-611							
	TPU:			+/-2200							
Cesium-137	12900			13000	pCi/L		101	(80%-120%)			
	Uncert:			+/-144							
	TPU:			+/-565							
Cobalt-60	10800			11300	pCi/L		104	(80%-120%)			
	Uncert:			+/-157							
	TPU:			+/-542							
Neptunium-237			U	42.2	pCi/L						
	Uncert:			+/-46.4							
	TPU:			+/-47.5							
Potassium-40			U	102	pCi/L						
	Uncert:			+/-114							
	TPU:			+/-116							
Sodium-22			U	-11	pCi/L						
	Uncert:			+/-15.7							
	TPU:			+/-15.9							
QC1203978021	MB										
Cesium-137			U	1.30	pCi/L				BSW1	02/23/18	15:41
	Uncert:			+/-1.31							
	TPU:			+/-1.34							
Cobalt-60			U	2.06	pCi/L						
	Uncert:			+/-0.951							

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Parmname	NOM	Sample	Qual	QC	Units	RER	REC%	Range	Anlst	Date	Time
Rad Gamma Spec											
Batch	1741844										
Neptunium-237	TPU:			+/-1.07							
			U	1.54	pCi/L						
	Uncert:			+/-1.99							
Potassium-40	TPU:			+/-2.03							
			U	-41	pCi/L						
	Uncert:			+/-15.1							
Sodium-22	TPU:			+/-17.9							
			U	0.391	pCi/L						
	Uncert:			+/-1.09							
	TPU:			+/-1.10							
Rad Gas Flow											
Batch	1742280										
QC1203978997	444396013	DUP									
Alpha	U	1.66	U	1.99	pCi/L	0.0957		(0-1)	BXG2	03/08/18	13:38
	Uncert:	+/-0.765		+/-0.906							
	TPU:	+/-0.778		+/-0.921							
Beta	U	1.06	U	1.10	pCi/L	0.0141		(0-1)		03/08/18	08:31
	Uncert:	+/-0.798		+/-0.720							
	TPU:	+/-0.804		+/-0.726							
QC1203979000	LCS										
Alpha	12.1			13.0	pCi/L		108	(80%-120%)	BXG2	03/08/18	13:41
	Uncert:			+/-0.626							
	TPU:			+/-1.30							
Beta	47.0			50.1	pCi/L		107	(80%-120%)		03/08/18	08:22
	Uncert:			+/-0.898							
	TPU:			+/-4.27							
QC1203978996	MB										
Alpha			U	-0.28	pCi/L				BXG2	03/08/18	13:38
	Uncert:			+/-0.0398							
	TPU:			+/-0.0399							
Beta			U	-0.0375	pCi/L					03/08/18	08:21
	Uncert:			+/-0.107							
	TPU:			+/-0.107							
QC1203978998	444396013	MS									
Alpha	483 U	1.66		442	pCi/L		91.4	(75%-125%)	BXG2	03/08/18	13:41
	Uncert:	+/-0.765		+/-22.4							
	TPU:	+/-0.778		+/-44.3							
Beta	1880 U	1.06		1970	pCi/L		105	(75%-125%)		03/08/18	08:21
	Uncert:	+/-0.798		+/-36.0							
	TPU:	+/-0.804		+/-171							
QC1203978999	444396013	MSD									
Alpha	483 U	1.66		473	pCi/L	0.172	97.9	(0-1)	BXG2	03/08/18	13:41
	Uncert:	+/-0.765		+/-24.1							
	TPU:	+/-0.778		+/-47.5							
Beta	1880 U	1.06		1910	pCi/L	0.0959	101	(0-1)		03/08/18	08:22
	Uncert:	+/-0.798		+/-34.4							

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Parmname	NOM	Sample	Qual	QC	Units	RER	REC%	Range	Anlst	Date	Time
Rad Gas Flow											
Batch	1742280										
		TPU:	+/-0.804	+/-161							
Batch	1742296										
QC1203979042	444617006	DUP									
Strontium-90		U	0.0107	U	-0.091	pCi/L	0.262	(0-1)	KSD1	02/28/18	16:16
		Uncert:	+/-0.118		+/-0.0763						
		TPU:	+/-0.118		+/-0.0763						
**Strontium Carrier		4.30	3.90		3.90	mg	90.7	(50%-105%)			
QC1203979044	LCS										
Strontium-90		23.5			23.3	pCi/L	99.1	(80%-120%)	KSD1	02/28/18	16:16
		Uncert:			+/-0.715						
		TPU:			+/-2.01						
**Strontium Carrier		4.30			3.30	mg	76.7	(50%-105%)			
QC1203979041	MB										
Strontium-90				U	-0.0129	pCi/L			KSD1	03/01/18	08:36
		Uncert:			+/-0.0885						
		TPU:			+/-0.0885						
**Strontium Carrier		4.30			4.10	mg	95.3	(50%-105%)			
QC1203979043	444617006	MS									
Strontium-90		235	U	0.0107	240	pCi/L	102	(75%-125%)	KSD1	02/28/18	16:16
		Uncert:		+/-0.118	+/-7.08						
		TPU:		+/-0.118	+/-20.5						
**Strontium Carrier		4.30	3.90		3.60	mg	83.7	(50%-105%)			

Notes:

TPU and Counting Uncertainty are calculated at the 68% confidence level (1-sigma).

The Qualifiers in this report are defined as follows:

- ** Analyte is a Tracer compound
- < Result is less than value reported
- > Result is greater than value reported
- BD Results are either below the MDC or tracer recovery is low
- FA Failed analysis.
- H Analytical holding time was exceeded
- J Value is estimated
- K Analyte present. Reported value may be biased high. Actual value is expected to be lower.
- L Analyte present. Reported value may be biased low. Actual value is expected to be higher.
- M M if above MDC and less than LLD
- M REMF Result > MDC/CL and < RDL
- N/A RPD or %Recovery limits do not apply.
- N1 See case narrative
- ND Analyte concentration is not detected above the detection limit
- NJ Consult Case Narrative, Data Summary package, or Project Manager concerning this qualifier
- Q One or more quality control criteria have not been met. Refer to the applicable narrative or DER.
- R Sample results are rejected

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Parmname	NOM	Sample Qual	QC	Units	RER	REC%	Range	Anlst	Date	Time
U	Analyte was analyzed for, but not detected above the MDL, MDA, MDC or LOD.									
UI	Gamma Spectroscopy--Uncertain identification									
UJ	Gamma Spectroscopy--Uncertain identification									
UL	Not considered detected. The associated number is the reported concentration, which may be inaccurate due to a low bias.									
X	Consult Case Narrative, Data Summary package, or Project Manager concerning this qualifier									
Y	Other specific qualifiers were required to properly define the results. Consult case narrative.									
^	RPD of sample and duplicate evaluated using +/-RL. Concentrations are <5X the RL. Qualifier Not Applicable for Radiochemistry.									
h	Preparation or preservation holding time was exceeded									

N/A indicates that spike recovery limits do not apply when sample concentration exceeds spike conc. by a factor of 4 or more or %RPD not applicable.

** Indicates analyte is a surrogate/tracer compound.

^ The Relative Percent Difference (RPD) obtained from the sample duplicate (DUP) is evaluated against the acceptance criteria when the sample is greater than five times (5X) the contract required detection limit (RL). In cases where either the sample or duplicate value is less than 5X the RL, a control limit of +/- the RL is used to evaluate the DUP result.

For PS, PSD, and SDILT results, the values listed are the measured amounts, not final concentrations.

Where the analytical method has been performed under NELAP certification, the analysis has met all of the requirements of the NELAC standard unless qualified on the QC Summary.