checked by N
6/21/17

CERTIFICATION

SDG No:	FA42192	Laboratory:	Accutest, Florida
Site:	BMS, Humacao, PR	Matrix:	AQ - Water

SUMMARY: Groundwater samples (Table 1) were collected on the BMSMC, Humacao, PR. Samples were taken March 16-17, 2017 and were analyzed in Accutest Laboratory of Orlando, Florida that reported the data under SDG No.: FA42192. Results were validated using the latest validation guidelines (July, 2015) of the EPA Hazardous Waste Support Section or the QC requirements of the method employed. The analyses performed are shown in Table 1. Individual data review worksheets are enclosed for each target analyte group. The organic data sample summary form shows for analyte results that were qualified.

In summary the results are valid and can be used for decision making purposes.

Table 1. Samples analyzed and analysis performed

SAMPLE ID	SAMPLE DESCRIPTION	MATRIX	ANALYSIS PERFORMED
FA42192-1	FB031617	AQ – Field Blank Water	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides
FA42192-2	S-34	Groundwater	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides
FA42192-3	S-33	Groundwater	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides
FA42192-3D	S-34 MSD	Groundwater	Pesticides
FA42192-3S	S-34 MS	Groundwater	Pesticides
FA42192-4	TB031617NR	AQ – Trip Blank Water	VOCs
FA42192-5	TB031617HR	AQ – Trip Blank Water	VOCs
FA42192-6	S-32	Groundwater	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides
FA42192-7	S-36	Groundwater	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides
FA42192-8	UP-1	Groundwater	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides
FA42192-9	TB031717NR	AQ – Trip Blank Water	VOCs
FA42192-10	TB031717HR	AQ – Trip Blank Water	VOCs
FA42192-11	UP-1DUP	Groundwater	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides
FA42192-12	S-28	Groundwater	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides

FA42192-13	UP-2	Groundwater	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides
FA42192-14	S-37	Groundwater	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides
FA42192-15	FB031717	AQ – Field Blank Water	VOCs; SVOCs; SVOCs (SIM); VPHs; EPHs; Pesticides

Reviewer Name: Rafael Infante
Chemist License 1888

Signature:

Date: April 26, 2017

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC. # 1888
QUIMICO LICENCIADO
$1.00
617222

SGS Accutest

Report of Analysis

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	Date Sampled:	03/16/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49212.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

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Rafael Infante Méndez
LIC. # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	Date Sampled:	03/16/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	15.1	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	99%		83-118%
17060-07-0	1,2-Dichloroethane-D4	99%		79-125%
2037-26-5	Toluene-D8	98%		85-112%
460-00-4	4-Bromofluorobenzene	97%		83-118%

(a) Associated ICV outside control limits.

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49213.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

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LIC. # 1008
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J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	4.1	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	31.6	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	98%		83-118%
17060-07-0	1,2-Dichloroethane-D4	97%		79-125%
2037-26-5	Toluene-D8	98%		85-112%
460-00-4	4-Bromofluorobenzene	99%		83-118%

(a) Associated ICV outside control limits.

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LIC. # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49214.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	0.45	1.0	0.20	ug/l	J
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

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ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	45.1	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	4.5	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	6.8	20	5.3	ug/l	J
75-65-0	Tert-Butyl Alcohol	33.3	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	99%		83-118%
17060-07-0	1,2-Dichloroethane-D4	99%		79-125%
2037-26-5	Toluene-D8	96%		85-112%
460-00-4	4-Bromofluorobenzene	98%		83-118%

(a) Associated ICV outside control limits.

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ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	TB031617NR		
Lab Sample ID:	FA42192-4	**Date Sampled:**	03/16/17
Matrix:	AQ - Trip Blank Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49215.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	TB031617NR		
Lab Sample ID:	FA42192-4	**Date Sampled:**	03/16/17
Matrix:	AQ - Trip Blank Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	1.8	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		83-118%
17060-07-0	1,2-Dichloroethane-D4	100%		79-125%
2037-26-5	Toluene-D8	97%		85-112%
460-00-4	4-Bromofluorobenzene	98%		83-118%

(a) Associated ICV outside control limits.

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC # 1888
QUIMICO LICENCIADO

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

SGS Accutest

Report of Analysis

Client Sample ID:	TB031617HR		
Lab Sample ID:	FA42192-5	Date Sampled:	03/16/17
Matrix:	AQ - Trip Blank Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49216.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC # 1888
QUIMICO LICENCIADO

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

Report of Analysis

Client Sample ID:	TB031617HR		
Lab Sample ID:	FA42192-5	**Date Sampled:**	03/16/17
Matrix:	AQ - Trip Blank Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	2.7	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		83-118%
17060-07-0	1,2-Dichloroethane-D4	99%		79-125%
2037-26-5	Toluene-D8	98%		85-112%
460-00-4	4-Bromofluorobenzene	98%		83-118%

(a) Associated ICV outside control limits.

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC # 1898
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49254.D	200	03/30/17	MM	n/a	n/a	VP1859
Run #2	P49217.D	2000	03/29/17	MM	n/a	n/a	VP1857

	Purge Volume
Run #1	5.0 ml
Run #2	5.0 ml

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	5000	2000	ug/l	
71-43-2	Benzene	ND	200	62	ug/l	
100-44-7	Benzyl Chloride	ND	400	73	ug/l	
74-97-5	Bromochloromethane	ND	200	90	ug/l	
75-27-4	Bromodichloromethane	ND	200	48	ug/l	
75-25-2	Bromoform	ND	200	81	ug/l	
106-99-0	1,3-Butadiene	ND	400	53	ug/l	
78-93-3	2-Butanone (MEK)	ND	1000	400	ug/l	
75-15-0	Carbon Disulfide	ND	400	110	ug/l	
56-23-5	Carbon Tetrachloride	ND	200	71	ug/l	
108-90-7	Chlorobenzene	ND	200	40	ug/l	
75-00-3	Chloroethane	ND	400	130	ug/l	
67-66-3	Chloroform	ND	200	60	ug/l	
110-82-7	Cyclohexane	ND	200	78	ug/l	
124-48-1	Dibromochloromethane	ND	200	55	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	1000	210	ug/l	
106-93-4	1,2-Dibromoethane	ND	400	55	ug/l	
75-71-8	Dichlorodifluoromethane	ND	400	100	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	200	65	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	200	43	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	200	51	ug/l	
75-34-3	1,1-Dichloroethane	ND	200	68	ug/l	
107-06-2	1,2-Dichloroethane	ND	200	62	ug/l	
75-35-4	1,1-Dichloroethylene	ND	200	64	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	200	55	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	200	44	ug/l	
78-87-5	1,2-Dichloropropane	ND	200	85	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	200	58	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	200	43	ug/l	
100-41-4	Ethylbenzene	34200 [a]	2000	710	ug/l	
76-13-1	Freon 113	ND	200	96	ug/l	
591-78-6	2-Hexanone	ND	2000	400	ug/l	

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
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E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	385	200	44	ug/l	
99-87-6	p-Isopropyltoluene	ND	200	43	ug/l	
79-20-9	Methyl Acetate	ND	4000	1000	ug/l	
74-83-9	Methyl Bromide	ND	400	120	ug/l	
74-87-3	Methyl Chloride	ND	400	100	ug/l	
108-87-2	Methylcyclohexane	ND	200	87	ug/l	
75-09-2	Methylene Chloride	976	1000	400	ug/l	J
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	1000	200	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	200	46	ug/l	
100-42-5	Styrene	ND	200	44	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	4000	1100	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	4000	1100	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	200	60	ug/l	
127-18-4	Tetrachloroethylene	ND	200	43	ug/l	
109-99-9	Tetrahydrofuran	ND	1000	320	ug/l	
108-88-3	Toluene	62.7	200	60	ug/l	J
87-61-6	1,2,3-Trichlorobenzene	ND	400	120	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	400	100	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	200	50	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	200	93	ug/l	
79-01-6	Trichloroethylene	ND	200	69	ug/l	
75-69-4	Trichlorofluoromethane	ND	400	100	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	200	65	ug/l	
75-01-4	Vinyl Chloride	ND	200	82	ug/l	
	m,p-Xylene	56300 [a]	4000	930	ug/l	
95-47-6	o-Xylene	3650 [a]	2000	510	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	99%	99%	83-118%
17060-07-0	1,2-Dichloroethane-D4	103%	101%	79-125%
2037-26-5	Toluene-D8	100%	99%	85-112%
460-00-4	4-Bromofluorobenzene	99%	100%	83-118%

(a) Result is from Run# 2

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC # 1888
QUIMICO LICENCIADO

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49218.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ESTADO LIBRE ASOCIADO DE PUERTO RICO • Rafael Infante Méndez LIC # 1886 • QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	8.0	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	0.61	1.0	0.23	ug/l	J
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	6.2	20	5.3	ug/l	J
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		83-118%
17060-07-0	1,2-Dichloroethane-D4	100%		79-125%
2037-26-5	Toluene-D8	97%		85-112%
460-00-4	4-Bromofluorobenzene	99%		83-118%

(a) Associated ICV outside control limits.

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC # 1666
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49219.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	36.9	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		83-118%
17060-07-0	1,2-Dichloroethane-D4	101%		79-125%
2037-26-5	Toluene-D8	95%		85-112%
460-00-4	4-Bromofluorobenzene	96%		83-118%

(a) Associated ICV outside control limits.

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	TB031717NR		
Lab Sample ID:	FA42192-9	Date Sampled:	03/17/17
Matrix:	AQ - Trip Blank Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49220.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	TB031717NR		
Lab Sample ID:	FA42192-9	Date Sampled:	03/17/17
Matrix:	AQ - Trip Blank Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	2.9	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	99%		83-118%
17060-07-0	1,2-Dichloroethane-D4	101%		79-125%
2037-26-5	Toluene-D8	98%		85-112%
460-00-4	4-Bromofluorobenzene	98%		83-118%

(a) Associated ICV outside control limits.

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1808
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit J = Indicates an estimated value
RL = Reporting Limit B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	TB031717HR		
Lab Sample ID:	FA42192-10	Date Sampled:	03/17/17
Matrix:	AQ - Trip Blank Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49221.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	TB031717HR		
Lab Sample ID:	FA42192-10	Date Sampled:	03/17/17
Matrix:	AQ - Trip Blank Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	3.1	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		83-118%
17060-07-0	1,2-Dichloroethane-D4	100%		79-125%
2037-26-5	Toluene-D8	98%		85-112%
460-00-4	4-Bromofluorobenzene	97%		83-118%

(a) Associated ICV outside control limits.

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49222.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	38.4	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	101%		83-118%
17060-07-0	1,2-Dichloroethane-D4	102%		79-125%
2037-26-5	Toluene-D8	94%		85-112%
460-00-4	4-Bromofluorobenzene	97%		83-118%

(a) Associated ICV outside control limits.

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	P49223.D	1	03/29/17	MM	n/a	n/a	VP1857
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane [a]	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropane	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ND = Not detected MDL = Method Detection Limit J = Indicates an estimated value
RL = Reporting Limit B = Indicates analyte found in associated method blank
E = Indicates value exceeds calibration range N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	1.4	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	100%		83-118%
17060-07-0	1,2-Dichloroethane-D4	101%		79-125%
2037-26-5	Toluene-D8	97%		85-112%
460-00-4	4-Bromofluorobenzene	98%		83-118%

(a) Associated ICV outside control limits.

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1[illegible]
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	M94807.D	1	03/29/17	SP	n/a	n/a	VM4067
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	0.46	1.0	0.20	ug/l	J
75-00-3	Chloroethane	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropan [a]	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	2.9	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	101%		83-118%
17060-07-0	1,2-Dichloroethane-D4	92%		79-125%
2037-26-5	Toluene-D8	91%		85-112%
460-00-4	4-Bromofluorobenzene	101%		83-118%

(a) Associated CCV outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	M94808.D	1	03/29/17	SP	n/a	n/a	VM4067
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropan [a]	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	0.74	1.0	0.23	ug/l	J
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	104%		83-118%
17060-07-0	1,2-Dichloroethane-D4	96%		79-125%
2037-26-5	Toluene-D8	92%		85-112%
460-00-4	4-Bromofluorobenzene	100%		83-118%

(a) Associated CCV outside control limits.

ISLAS DEL LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
LIC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	Date Sampled:	03/17/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846 8260C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	M94806.D	1	03/29/17	SP	n/a	n/a	VM4067
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
67-64-1	Acetone	ND	25	10	ug/l	
71-43-2	Benzene	ND	1.0	0.31	ug/l	
100-44-7	Benzyl Chloride	ND	2.0	0.36	ug/l	
74-97-5	Bromochloromethane	ND	1.0	0.45	ug/l	
75-27-4	Bromodichloromethane	ND	1.0	0.24	ug/l	
75-25-2	Bromoform	ND	1.0	0.41	ug/l	
106-99-0	1,3-Butadiene	ND	2.0	0.26	ug/l	
78-93-3	2-Butanone (MEK)	ND	5.0	2.0	ug/l	
75-15-0	Carbon Disulfide	ND	2.0	0.53	ug/l	
56-23-5	Carbon Tetrachloride	ND	1.0	0.36	ug/l	
108-90-7	Chlorobenzene	ND	1.0	0.20	ug/l	
75-00-3	Chloroethane	ND	2.0	0.67	ug/l	
67-66-3	Chloroform	ND	1.0	0.30	ug/l	
110-82-7	Cyclohexane	ND	1.0	0.39	ug/l	
124-48-1	Dibromochloromethane	ND	1.0	0.28	ug/l	
96-12-8	1,2-Dibromo-3-chloropropan [a]	ND	5.0	1.0	ug/l	
106-93-4	1,2-Dibromoethane	ND	2.0	0.28	ug/l	
75-71-8	Dichlorodifluoromethane	ND	2.0	0.50	ug/l	
95-50-1	1,2-Dichlorobenzene	ND	1.0	0.32	ug/l	
541-73-1	1,3-Dichlorobenzene	ND	1.0	0.22	ug/l	
106-46-7	1,4-Dichlorobenzene	ND	1.0	0.26	ug/l	
75-34-3	1,1-Dichloroethane	ND	1.0	0.34	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.31	ug/l	
75-35-4	1,1-Dichloroethylene	ND	1.0	0.32	ug/l	
156-59-2	cis-1,2-Dichloroethylene	ND	1.0	0.28	ug/l	
156-60-5	trans-1,2-Dichloroethylene	ND	1.0	0.22	ug/l	
78-87-5	1,2-Dichloropropane	ND	1.0	0.43	ug/l	
10061-01-5	cis-1,3-Dichloropropene	ND	1.0	0.29	ug/l	
10061-02-6	trans-1,3-Dichloropropene	ND	1.0	0.21	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.36	ug/l	
76-13-1	Freon 113	ND	1.0	0.48	ug/l	
591-78-6	2-Hexanone	ND	10	2.0	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1888
QUIMICO LICENCIADO

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	**Date Sampled:**	03/17/17
Matrix:	AQ - Field Blank Water	**Date Received:**	03/18/17
Method:	SW846 8260C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

VOA TCL List (SOM02.0)

CAS No.	Compound	Result	RL	MDL	Units	Q
98-82-8	Isopropylbenzene	ND	1.0	0.22	ug/l	
99-87-6	p-Isopropyltoluene	ND	1.0	0.21	ug/l	
79-20-9	Methyl Acetate	ND	20	5.0	ug/l	
74-83-9	Methyl Bromide	ND	2.0	0.59	ug/l	
74-87-3	Methyl Chloride	ND	2.0	0.50	ug/l	
108-87-2	Methylcyclohexane	ND	1.0	0.44	ug/l	
75-09-2	Methylene Chloride	ND	5.0	2.0	ug/l	
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0	1.0	ug/l	
1634-04-4	Methyl Tert Butyl Ether	ND	1.0	0.23	ug/l	
100-42-5	Styrene	ND	1.0	0.22	ug/l	
75-85-4	Tert-Amyl Alcohol	ND	20	5.3	ug/l	
75-65-0	Tert-Butyl Alcohol	ND	20	5.3	ug/l	
79-34-5	1,1,2,2-Tetrachloroethane	ND	1.0	0.30	ug/l	
127-18-4	Tetrachloroethylene	ND	1.0	0.22	ug/l	
109-99-9	Tetrahydrofuran	ND	5.0	1.6	ug/l	
108-88-3	Toluene	ND	1.0	0.30	ug/l	
87-61-6	1,2,3-Trichlorobenzene	ND	2.0	0.61	ug/l	
120-82-1	1,2,4-Trichlorobenzene	ND	2.0	0.50	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.47	ug/l	
79-01-6	Trichloroethylene	ND	1.0	0.35	ug/l	
75-69-4	Trichlorofluoromethane	ND	2.0	0.50	ug/l	
95-63-6	1,2,4-Trimethylbenzene	ND	1.0	0.32	ug/l	
75-01-4	Vinyl Chloride	ND	1.0	0.41	ug/l	
	m,p-Xylene	ND	2.0	0.47	ug/l	
95-47-6	o-Xylene	ND	1.0	0.26	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	102%		83-118%
17060-07-0	1,2-Dichloroethane-D4	91%		79-125%
2037-26-5	Toluene-D8	93%		85-112%
460-00-4	4-Bromofluorobenzene	98%		83-118%

(a) Associated CCV outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	Date Sampled:	03/16/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690711.D	1	03/30/17	NG	03/23/17	OP64302	SL3980
Run #2							

	Initial Volume	Final Volume
Run #1	1020 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	9.8	2.5	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	4.9	0.58	ug/l	
95-57-8	2-Chlorophenol	ND	4.9	0.62	ug/l	
120-83-2	2,4-Dichlorophenol	ND	4.9	0.82	ug/l	
105-67-9	2,4-Dimethylphenol	ND	4.9	0.72	ug/l	
51-28-5	2,4-Dinitrophenol [a]	ND	25	4.9	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	9.8	2.0	ug/l	
95-48-7	2-Methylphenol	ND	4.9	0.55	ug/l	
	3&4-Methylphenol	ND	4.9	0.96	ug/l	
88-75-5	2-Nitrophenol [a]	ND	4.9	0.84	ug/l	
100-02-7	4-Nitrophenol	ND	25	4.9	ug/l	
87-86-5	Pentachlorophenol	ND	25	4.9	ug/l	
108-95-2	Phenol	ND	4.9	0.49	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol [a]	ND	4.9	0.95	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	4.9	0.73	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	4.9	0.74	ug/l	
83-32-9	Acenaphthene	ND	4.9	0.61	ug/l	
208-96-8	Acenaphthylene	ND	4.9	0.63	ug/l	
98-86-2	Acetophenone	ND	4.9	0.79	ug/l	
120-12-7	Anthracene	ND	4.9	0.78	ug/l	
1912-24-9	Atrazine	ND	4.9	1.0	ug/l	
100-52-7	Benzaldehyde	ND	25	4.9	ug/l	
56-55-3	Benzo(a)anthracene	ND	4.9	0.75	ug/l	
50-32-8	Benzo(a)pyrene	ND	4.9	0.77	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	4.9	0.76	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	4.9	0.81	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	4.9	0.84	ug/l	
92-52-4	1,1'-Biphenyl	ND	4.9	0.61	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	4.9	0.83	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	4.9	0.98	ug/l	
86-74-8	Carbazole	ND	4.9	0.59	ug/l	
106-47-8	4-Chloroaniline	ND	4.9	0.62	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IQ # 1888
QUIMICO LICENCIADO

Report of Analysis

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	**Date Sampled:**	03/16/17
Matrix:	AQ - Field Blank Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	4.9	0.79	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	4.9	0.72	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	4.9	0.74	ug/l	
91-58-7	2-Chloronaphthalene	ND	4.9	0.49	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	4.9	0.53	ug/l	
218-01-9	Chrysene	ND	4.9	0.83	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	4.9	0.79	ug/l	
132-64-9	Dibenzofuran	ND	4.9	0.59	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	4.9	0.63	ug/l	
84-66-2	Diethyl Phthalate	ND	4.9	0.98	ug/l	
131-11-3	Dimethyl Phthalate	ND	4.9	0.98	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	4.9	0.98	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	4.9	0.98	ug/l	
121-14-2	2,4-Dinitrotoluene [a]	ND	4.9	0.80	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	4.9	0.70	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	4.9	0.98	ug/l	
206-44-0	Fluoranthene	ND	4.9	0.54	ug/l	
86-73-7	Fluorene	ND	4.9	0.69	ug/l	
118-74-1	Hexachlorobenzene	ND	4.9	0.68	ug/l	
87-68-3	Hexachlorobutadiene	ND	4.9	0.49	ug/l	
77-47-4	Hexachlorocyclopentadiene [a]	ND	4.9	1.8	ug/l	
67-72-1	Hexachloroethane	ND	4.9	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	4.9	0.70	ug/l	
78-59-1	Isophorone	ND	4.9	0.76	ug/l	
91-57-6	2-Methylnaphthalene	ND	4.9	0.59	ug/l	
88-74-4	2-Nitroaniline [a]	ND	4.9	1.8	ug/l	
99-09-2	3-Nitroaniline [a]	ND	4.9	0.86	ug/l	
100-01-6	4-Nitroaniline	ND	4.9	1.1	ug/l	
98-95-3	Nitrobenzene	ND	4.9	0.91	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	4.9	0.66	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	4.9	0.79	ug/l	
85-01-8	Phenanthrene	ND	4.9	0.85	ug/l	
129-00-0	Pyrene	ND	4.9	0.67	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	4.9	0.49	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	24%		14-67%
4165-62-2	Phenol-d5	17%		10-50%
118-79-6	2,4,6-Tribromophenol	92%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	**Date Sampled:**	03/16/17
Matrix:	AQ - Field Blank Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	75%		42-108%
321-60-8	2-Fluorobiphenyl	63%		40-106%
1718-51-0	Terphenyl-d14	65%		39-121%

(a) Associated CCV outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690712.D	1	03/30/17	NG	03/23/17	OP64302	SL3980
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	10	2.5	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	5.0	0.59	ug/l	
95-57-8	2-Chlorophenol	ND	5.0	0.63	ug/l	
120-83-2	2,4-Dichlorophenol	ND	5.0	0.84	ug/l	
105-67-9	2,4-Dimethylphenol	ND	5.0	0.74	ug/l	
51-28-5	2,4-Dinitrophenol [a]	ND	25	5.0	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	10	2.0	ug/l	
95-48-7	2-Methylphenol	ND	5.0	0.56	ug/l	
	3&4-Methylphenol	ND	5.0	0.98	ug/l	
88-75-5	2-Nitrophenol [a]	ND	5.0	0.85	ug/l	
100-02-7	4-Nitrophenol	ND	25	5.0	ug/l	
87-86-5	Pentachlorophenol	ND	25	5.0	ug/l	
108-95-2	Phenol	ND	5.0	0.50	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol [a]	ND	5.0	0.97	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.74	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.75	ug/l	
83-32-9	Acenaphthene	ND	5.0	0.63	ug/l	
208-96-8	Acenaphthylene	ND	5.0	0.64	ug/l	
98-86-2	Acetophenone	ND	5.0	0.81	ug/l	
120-12-7	Anthracene	ND	5.0	0.80	ug/l	
1912-24-9	Atrazine	ND	5.0	1.0	ug/l	
100-52-7	Benzaldehyde	ND	25	5.0	ug/l	
56-55-3	Benzo(a)anthracene	ND	5.0	0.76	ug/l	
50-32-8	Benzo(a)pyrene	ND	5.0	0.78	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	5.0	0.78	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	5.0	0.82	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	5.0	0.86	ug/l	
92-52-4	1,1'-Biphenyl	ND	5.0	0.62	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	5.0	0.85	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	5.0	1.0	ug/l	
86-74-8	Carbazole	ND	5.0	0.60	ug/l	
106-47-8	4-Chloroaniline	ND	5.0	0.63	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	5.0	0.81	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	5.0	0.73	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	5.0	0.76	ug/l	
91-58-7	2-Chloronaphthalene	ND	5.0	0.50	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	5.0	0.54	ug/l	
218-01-9	Chrysene	ND	5.0	0.85	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	5.0	0.80	ug/l	
132-64-9	Dibenzofuran	ND	5.0	0.60	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	5.0	0.64	ug/l	
84-66-2	Diethyl Phthalate	ND	5.0	1.0	ug/l	
131-11-3	Dimethyl Phthalate	ND	5.0	1.0	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	5.0	1.0	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	5.0	1.0	ug/l	
121-14-2	2,4-Dinitrotoluene [a]	ND	5.0	0.81	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	5.0	0.71	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	5.0	1.0	ug/l	
206-44-0	Fluoranthene	ND	5.0	0.55	ug/l	
86-73-7	Fluorene	ND	5.0	0.70	ug/l	
118-74-1	Hexachlorobenzene	ND	5.0	0.69	ug/l	
87-68-3	Hexachlorobutadiene	ND	5.0	0.50	ug/l	
77-47-4	Hexachlorocyclopentadiene [a]	ND	5.0	1.8	ug/l	
67-72-1	Hexachloroethane	ND	5.0	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	5.0	0.71	ug/l	
78-59-1	Isophorone	ND	5.0	0.78	ug/l	
91-57-6	2-Methylnaphthalene	ND	5.0	0.60	ug/l	
88-74-4	2-Nitroaniline [a]	ND	5.0	1.8	ug/l	
99-09-2	3-Nitroaniline [a]	ND	5.0	0.88	ug/l	
100-01-6	4-Nitroaniline	ND	5.0	1.2	ug/l	
98-95-3	Nitrobenzene	ND	5.0	0.93	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	5.0	0.67	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	5.0	0.81	ug/l	
85-01-8	Phenanthrene	ND	5.0	0.86	ug/l	
129-00-0	Pyrene	ND	5.0	0.68	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	5.0	0.50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	29%		14-67%
4165-62-2	Phenol-d5	22%		10-50%
118-79-6	2,4,6-Tribromophenol	109%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	84%		42-108%
321-60-8	2-Fluorobiphenyl	73%		40-106%
1718-51-0	Terphenyl-d14	66%		39-121%

(a) Associated CCV outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690713.D	1	03/30/17	NG	03/23/17	OP64302	SL3980
Run #2							

	Initial Volume	Final Volume
Run #1	1020 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	9.8	2.5	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	4.9	0.58	ug/l	
95-57-8	2-Chlorophenol	ND	4.9	0.62	ug/l	
120-83-2	2,4-Dichlorophenol	ND	4.9	0.82	ug/l	
105-67-9	2,4-Dimethylphenol	ND	4.9	0.72	ug/l	
51-28-5	2,4-Dinitrophenol [a]	ND	25	4.9	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	9.8	2.0	ug/l	
95-48-7	2-Methylphenol	ND	4.9	0.55	ug/l	
	3&4-Methylphenol	ND	4.9	0.96	ug/l	
88-75-5	2-Nitrophenol [a]	ND	4.9	0.84	ug/l	
100-02-7	4-Nitrophenol	ND	25	4.9	ug/l	
87-86-5	Pentachlorophenol	ND	25	4.9	ug/l	
108-95-2	Phenol	ND	4.9	0.49	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol [a]	ND	4.9	0.95	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	4.9	0.73	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	4.9	0.74	ug/l	
83-32-9	Acenaphthene	ND	4.9	0.61	ug/l	
208-96-8	Acenaphthylene	ND	4.9	0.63	ug/l	
98-86-2	Acetophenone	ND	4.9	0.79	ug/l	
120-12-7	Anthracene	ND	4.9	0.78	ug/l	
1912-24-9	Atrazine	ND	4.9	1.0	ug/l	
100-52-7	Benzaldehyde	ND	25	4.9	ug/l	
56-55-3	Benzo(a)anthracene	ND	4.9	0.75	ug/l	
50-32-8	Benzo(a)pyrene	ND	4.9	0.77	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	4.9	0.76	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	4.9	0.81	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	4.9	0.84	ug/l	
92-52-4	1,1'-Biphenyl	ND	4.9	0.61	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	4.9	0.83	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	4.9	0.98	ug/l	
86-74-8	Carbazole	ND	4.9	0.59	ug/l	
106-47-8	4-Chloroaniline	ND	4.9	0.62	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IQ # 1888
QUIMICO LICENCIADO

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	4.9	0.79	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	4.9	0.72	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	4.9	0.74	ug/l	
91-58-7	2-Chloronaphthalene	ND	4.9	0.49	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	4.9	0.53	ug/l	
218-01-9	Chrysene	ND	4.9	0.83	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	4.9	0.79	ug/l	
132-64-9	Dibenzofuran	ND	4.9	0.59	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	4.9	0.63	ug/l	
84-66-2	Diethyl Phthalate	ND	4.9	0.98	ug/l	
131-11-3	Dimethyl Phthalate	ND	4.9	0.98	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	4.9	0.98	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	4.9	0.98	ug/l	
121-14-2	2,4-Dinitrotoluene [a]	ND	4.9	0.80	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	4.9	0.70	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	4.9	0.98	ug/l	
206-44-0	Fluoranthene	ND	4.9	0.54	ug/l	
86-73-7	Fluorene	ND	4.9	0.69	ug/l	
118-74-1	Hexachlorobenzene	ND	4.9	0.68	ug/l	
87-68-3	Hexachlorobutadiene	ND	4.9	0.49	ug/l	
77-47-4	Hexachlorocyclopentadiene [a]	ND	4.9	1.8	ug/l	
67-72-1	Hexachloroethane	ND	4.9	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	4.9	0.70	ug/l	
78-59-1	Isophorone	ND	4.9	0.76	ug/l	
91-57-6	2-Methylnaphthalene	ND	4.9	0.59	ug/l	
88-74-4	2-Nitroaniline [a]	ND	4.9	1.8	ug/l	
99-09-2	3-Nitroaniline [a]	ND	4.9	0.86	ug/l	
100-01-6	4-Nitroaniline	ND	4.9	1.1	ug/l	
98-95-3	Nitrobenzene	ND	4.9	0.91	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	4.9	0.66	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	4.9	0.79	ug/l	
85-01-8	Phenanthrene	1.5	4.9	0.85	ug/l	J
129-00-0	Pyrene	ND	4.9	0.67	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	4.9	0.49	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	34%		14-67%
4165-62-2	Phenol-d5	20%		10-50%
118-79-6	2,4,6-Tribromophenol	103%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	86%		42-108%
321-60-8	2-Fluorobiphenyl	75%		40-106%
1718-51-0	Terphenyl-d14	67%		39-121%

(a) Associated CCV outside control limits.

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	L0690726.D	10	03/31/17	NG	03/23/17	OP64302	SL3981
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	100	25	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	50	5.9	ug/l	
95-57-8	2-Chlorophenol	ND	50	6.3	ug/l	
120-83-2	2,4-Dichlorophenol	ND	50	8.4	ug/l	
105-67-9	2,4-Dimethylphenol	53.2	50	7.4	ug/l	
51-28-5	2,4-Dinitrophenol [b]	ND	250	50	ug/l	
534-52-1	4,6-Dinitro-o-cresol [c]	ND	100	20	ug/l	
95-48-7	2-Methylphenol	ND	50	5.6	ug/l	
	3&4-Methylphenol	ND	50	9.8	ug/l	
88-75-5	2-Nitrophenol [b]	ND	50	8.5	ug/l	
100-02-7	4-Nitrophenol	ND	250	50	ug/l	
87-86-5	Pentachlorophenol [d]	ND	250	50	ug/l	
108-95-2	Phenol	ND	50	5.0	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol [d]	ND	50	9.7	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	50	7.4	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	50	7.5	ug/l	
83-32-9	Acenaphthene	ND	50	6.3	ug/l	
208-96-8	Acenaphthylene	ND	50	6.4	ug/l	
98-86-2	Acetophenone	27.9	50	8.1	ug/l	J
120-12-7	Anthracene	ND	50	8.0	ug/l	
1912-24-9	Atrazine	ND	50	10	ug/l	
100-52-7	Benzaldehyde	ND	250	50	ug/l	
56-55-3	Benzo(a)anthracene	ND	50	7.6	ug/l	
50-32-8	Benzo(a)pyrene	ND	50	7.8	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	50	7.8	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	50	8.2	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	50	8.6	ug/l	
92-52-4	1,1'-Biphenyl	ND	50	6.2	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	50	8.5	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	50	10	ug/l	
86-74-8	Carbazole	ND	50	6.0	ug/l	
106-47-8	4-Chloroaniline	ND	50	6.3	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	50	8.1	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	50	7.3	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	50	7.6	ug/l	
91-58-7	2-Chloronaphthalene	ND	50	5.0	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	50	5.4	ug/l	
218-01-9	Chrysene	ND	50	8.5	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	50	8.0	ug/l	
132-64-9	Dibenzofuran	ND	50	6.0	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	50	6.4	ug/l	
84-66-2	Diethyl Phthalate	ND	50	10	ug/l	
131-11-3	Dimethyl Phthalate	ND	50	10	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	50	10	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	50	10	ug/l	
121-14-2	2,4-Dinitrotoluene [b]	ND	50	8.1	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	50	7.1	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	50	10	ug/l	
206-44-0	Fluoranthene	ND	50	5.5	ug/l	
86-73-7	Fluorene	ND	50	7.0	ug/l	
118-74-1	Hexachlorobenzene	ND	50	6.9	ug/l	
87-68-3	Hexachlorobutadiene	ND	50	5.0	ug/l	
77-47-4	Hexachlorocyclopentadiene [d]	ND	50	18	ug/l	
67-72-1	Hexachloroethane	ND	50	16	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	50	7.1	ug/l	
78-59-1	Isophorone	ND	50	7.8	ug/l	
91-57-6	2-Methylnaphthalene	ND	50	6.0	ug/l	
88-74-4	2-Nitroaniline [d]	ND	50	18	ug/l	
99-09-2	3-Nitroaniline [c]	ND	50	8.8	ug/l	
100-01-6	4-Nitroaniline	ND	50	12	ug/l	
98-95-3	Nitrobenzene	ND	50	9.3	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	50	6.7	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	50	8.1	ug/l	
85-01-8	Phenanthrene	ND	50	8.6	ug/l	
129-00-0	Pyrene	ND	50	6.8	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	50	5.0	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	5% [e]		14-67%
4165-62-2	Phenol-d5	23%		10-50%
118-79-6	2,4,6-Tribromophenol	122% [e]		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	108%		42-108%
321-60-8	2-Fluorobiphenyl	83%		40-106%
1718-51-0	Terphenyl-d14	85%		39-121%

(a) Dilution required due to matrix interference.
(b) Associated CCV and BS outside control limits.
(c) Associated BS recovery outside control limits.
(d) Associated CCV outside control limits.
(e) Outside control limits due to matrix interference.

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Manuel Infante Méndez
IQ # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690717.D	1	03/31/17	NG	03/23/17	OP64302	SL3980
Run #2							

	Initial Volume	Final Volume
Run #1	1040 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	9.6	2.4	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	4.8	0.57	ug/l	
95-57-8	2-Chlorophenol	ND	4.8	0.61	ug/l	
120-83-2	2,4-Dichlorophenol	ND	4.8	0.80	ug/l	
105-67-9	2,4-Dimethylphenol	ND	4.8	0.71	ug/l	
51-28-5	2,4-Dinitrophenol [a]	ND	24	4.8	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	9.6	1.9	ug/l	
95-48-7	2-Methylphenol	ND	4.8	0.54	ug/l	
	3&4-Methylphenol	ND	4.8	0.94	ug/l	
88-75-5	2-Nitrophenol [a]	ND	4.8	0.82	ug/l	
100-02-7	4-Nitrophenol	ND	24	4.8	ug/l	
87-86-5	Pentachlorophenol	ND	24	4.8	ug/l	
108-95-2	Phenol	ND	4.8	0.48	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol [a]	ND	4.8	0.93	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	4.8	0.71	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	4.8	0.72	ug/l	
83-32-9	Acenaphthene	ND	4.8	0.60	ug/l	
208-96-8	Acenaphthylene	ND	4.8	0.61	ug/l	
98-86-2	Acetophenone	ND	4.8	0.78	ug/l	
120-12-7	Anthracene	ND	4.8	0.77	ug/l	
1912-24-9	Atrazine	ND	4.8	0.98	ug/l	
100-52-7	Benzaldehyde	ND	24	4.8	ug/l	
56-55-3	Benzo(a)anthracene	ND	4.8	0.73	ug/l	
50-32-8	Benzo(a)pyrene	ND	4.8	0.75	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	4.8	0.75	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	4.8	0.79	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	4.8	0.82	ug/l	
92-52-4	1,1'-Biphenyl	ND	4.8	0.60	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	4.8	0.81	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	4.8	0.96	ug/l	
86-74-8	Carbazole	ND	4.8	0.58	ug/l	
106-47-8	4-Chloroaniline	ND	4.8	0.61	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
1888
QUIMICO LICENCIADO

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	4.8	0.78	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	4.8	0.70	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	4.8	0.73	ug/l	
91-58-7	2-Chloronaphthalene	ND	4.8	0.48	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	4.8	0.52	ug/l	
218-01-9	Chrysene	ND	4.8	0.82	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	4.8	0.77	ug/l	
132-64-9	Dibenzofuran	ND	4.8	0.58	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	4.8	0.62	ug/l	
84-66-2	Diethyl Phthalate	ND	4.8	0.96	ug/l	
131-11-3	Dimethyl Phthalate	ND	4.8	0.96	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	4.8	0.96	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	4.8	0.96	ug/l	
121-14-2	2,4-Dinitrotoluene [a]	ND	4.8	0.78	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	4.8	0.69	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	4.8	0.96	ug/l	
206-44-0	Fluoranthene	ND	4.8	0.53	ug/l	
86-73-7	Fluorene	ND	4.8	0.67	ug/l	
118-74-1	Hexachlorobenzene	ND	4.8	0.67	ug/l	
87-68-3	Hexachlorobutadiene	ND	4.8	0.48	ug/l	
77-47-4	Hexachlorocyclopentadiene [a]	ND	4.8	1.7	ug/l	
67-72-1	Hexachloroethane	ND	4.8	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	4.8	0.69	ug/l	
78-59-1	Isophorone	ND	4.8	0.75	ug/l	
91-57-6	2-Methylnaphthalene	ND	4.8	0.58	ug/l	
88-74-4	2-Nitroaniline [a]	ND	4.8	1.7	ug/l	
99-09-2	3-Nitroaniline [a]	ND	4.8	0.85	ug/l	
100-01-6	4-Nitroaniline	ND	4.8	1.1	ug/l	
98-95-3	Nitrobenzene	ND	4.8	0.90	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	4.8	0.64	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	4.8	0.78	ug/l	
85-01-8	Phenanthrene	ND	4.8	0.83	ug/l	
129-00-0	Pyrene	ND	4.8	0.66	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	4.8	0.48	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	27%		14-67%
4165-62-2	Phenol-d5	19%		10-50%
118-79-6	2,4,6-Tribromophenol	107%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

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ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	86%		42-108%
321-60-8	2-Fluorobiphenyl	73%		40-106%
1718-51-0	Terphenyl-d14	69%		39-121%

(a) Associated CCV outside control limits.

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IQ # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
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E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690594.D	1	03/26/17	NG	03/24/17	OP64319	SL3976
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	10	2.5	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	5.0	0.59	ug/l	
95-57-8	2-Chlorophenol	ND	5.0	0.63	ug/l	
120-83-2	2,4-Dichlorophenol	ND	5.0	0.84	ug/l	
105-67-9	2,4-Dimethylphenol	ND	5.0	0.74	ug/l	
51-28-5	2,4-Dinitrophenol	ND	25	5.0	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	10	2.0	ug/l	
95-48-7	2-Methylphenol	ND	5.0	0.56	ug/l	
	3&4-Methylphenol	ND	5.0	0.98	ug/l	
88-75-5	2-Nitrophenol	ND	5.0	0.85	ug/l	
100-02-7	4-Nitrophenol	ND	25	5.0	ug/l	
87-86-5	Pentachlorophenol	ND	25	5.0	ug/l	
108-95-2	Phenol	ND	5.0	0.50	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.97	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.74	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.75	ug/l	
83-32-9	Acenaphthene	ND	5.0	0.63	ug/l	
208-96-8	Acenaphthylene	ND	5.0	0.64	ug/l	
98-86-2	Acetophenone	ND	5.0	0.81	ug/l	
120-12-7	Anthracene	1.3	5.0	0.80	ug/l	J
1912-24-9	Atrazine	ND	5.0	1.0	ug/l	
100-52-7	Benzaldehyde	ND	25	5.0	ug/l	
56-55-3	Benzo(a)anthracene	ND	5.0	0.76	ug/l	
50-32-8	Benzo(a)pyrene	ND	5.0	0.78	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	5.0	0.78	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	5.0	0.82	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	5.0	0.86	ug/l	
92-52-4	1,1'-Biphenyl	ND	5.0	0.62	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	5.0	0.85	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	5.0	1.0	ug/l	
86-74-8	Carbazole	ND	5.0	0.60	ug/l	
106-47-8	4-Chloroaniline	ND	5.0	0.63	ug/l	

ND = Not detected MDL = Method Detection Limit
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E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	5.0	0.81	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	5.0	0.73	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	5.0	0.76	ug/l	
91-58-7	2-Chloronaphthalene	ND	5.0	0.50	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	5.0	0.54	ug/l	
218-01-9	Chrysene	ND	5.0	0.85	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	5.0	0.80	ug/l	
132-64-9	Dibenzofuran	ND	5.0	0.60	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	5.0	0.64	ug/l	
84-66-2	Diethyl Phthalate	ND	5.0	1.0	ug/l	
131-11-3	Dimethyl Phthalate	ND	5.0	1.0	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	5.0	1.0	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	5.0	1.0	ug/l	
121-14-2	2,4-Dinitrotoluene	ND	5.0	0.81	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	5.0	0.71	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	5.0	1.0	ug/l	
206-44-0	Fluoranthene	ND	5.0	0.55	ug/l	
86-73-7	Fluorene	ND	5.0	0.70	ug/l	
118-74-1	Hexachlorobenzene	ND	5.0	0.69	ug/l	
87-68-3	Hexachlorobutadiene	ND	5.0	0.50	ug/l	
77-47-4	Hexachlorocyclopentadiene	ND	5.0	1.8	ug/l	
67-72-1	Hexachloroethane	ND	5.0	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	5.0	0.71	ug/l	
78-59-1	Isophorone	ND	5.0	0.78	ug/l	
91-57-6	2-Methylnaphthalene	ND	5.0	0.60	ug/l	
88-74-4	2-Nitroaniline	ND	5.0	1.8	ug/l	
99-09-2	3-Nitroaniline	ND	5.0	0.88	ug/l	
100-01-6	4-Nitroaniline	ND	5.0	1.2	ug/l	
98-95-3	Nitrobenzene	ND	5.0	0.93	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	5.0	0.67	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	5.0	0.81	ug/l	
85-01-8	Phenanthrene	ND	5.0	0.86	ug/l	
129-00-0	Pyrene	ND	5.0	0.68	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	5.0	0.50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	23%		14-67%
4165-62-2	Phenol-d5	14%		10-50%
118-79-6	2,4,6-Tribromophenol	71%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	58%		42-108%
321-60-8	2-Fluorobiphenyl	60%		40-106%
1718-51-0	Terphenyl-d14	57%		39-121%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
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J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690595.D	1	03/26/17	NG	03/24/17	OP64319	SL3976
Run #2							

	Initial Volume	Final Volume
Run #1	1020 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	9.8	2.5	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	4.9	0.58	ug/l	
95-57-8	2-Chlorophenol	ND	4.9	0.62	ug/l	
120-83-2	2,4-Dichlorophenol	ND	4.9	0.82	ug/l	
105-67-9	2,4-Dimethylphenol	ND	4.9	0.72	ug/l	
51-28-5	2,4-Dinitrophenol	ND	25	4.9	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	9.8	2.0	ug/l	
95-48-7	2-Methylphenol	ND	4.9	0.55	ug/l	
	3&4-Methylphenol	ND	4.9	0.96	ug/l	
88-75-5	2-Nitrophenol	ND	4.9	0.84	ug/l	
100-02-7	4-Nitrophenol	ND	25	4.9	ug/l	
87-86-5	Pentachlorophenol	ND	25	4.9	ug/l	
108-95-2	Phenol	ND	4.9	0.49	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol	ND	4.9	0.95	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	4.9	0.73	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	4.9	0.74	ug/l	
83-32-9	Acenaphthene	ND	4.9	0.61	ug/l	
208-96-8	Acenaphthylene	ND	4.9	0.63	ug/l	
98-86-2	Acetophenone	ND	4.9	0.79	ug/l	
120-12-7	Anthracene	1.4	4.9	0.78	ug/l	J
1912-24-9	Atrazine	ND	4.9	1.0	ug/l	
100-52-7	Benzaldehyde	ND	25	4.9	ug/l	
56-55-3	Benzo(a)anthracene	ND	4.9	0.75	ug/l	
50-32-8	Benzo(a)pyrene	ND	4.9	0.77	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	4.9	0.76	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	4.9	0.81	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	4.9	0.84	ug/l	
92-52-4	1,1'-Biphenyl	ND	4.9	0.61	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	4.9	0.83	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	4.9	0.98	ug/l	
86-74-8	Carbazole	ND	4.9	0.59	ug/l	
106-47-8	4-Chloroaniline	ND	4.9	0.62	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

LIBRE ASOCIACIÓN DE PUERTO RICO
Rafael Infante Méndez
IQ # 1888
QUÍMICO LICENCIADO

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	4.9	0.79	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	4.9	0.72	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	4.9	0.74	ug/l	
91-58-7	2-Chloronaphthalene	ND	4.9	0.49	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	4.9	0.53	ug/l	
218-01-9	Chrysene	ND	4.9	0.83	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	4.9	0.79	ug/l	
132-64-9	Dibenzofuran	ND	4.9	0.59	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	4.9	0.63	ug/l	
84-66-2	Diethyl Phthalate	ND	4.9	0.98	ug/l	
131-11-3	Dimethyl Phthalate	ND	4.9	0.98	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	4.9	0.98	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	4.9	0.98	ug/l	
121-14-2	2,4-Dinitrotoluene	ND	4.9	0.80	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	4.9	0.70	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	4.9	0.98	ug/l	
206-44-0	Fluoranthene	ND	4.9	0.54	ug/l	
86-73-7	Fluorene	ND	4.9	0.69	ug/l	
118-74-1	Hexachlorobenzene	ND	4.9	0.68	ug/l	
87-68-3	Hexachlorobutadiene	ND	4.9	0.49	ug/l	
77-47-4	Hexachlorocyclopentadiene	ND	4.9	1.8	ug/l	
67-72-1	Hexachloroethane	ND	4.9	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	4.9	0.70	ug/l	
78-59-1	Isophorone	ND	4.9	0.76	ug/l	
91-57-6	2-Methylnaphthalene	ND	4.9	0.59	ug/l	
88-74-4	2-Nitroaniline	ND	4.9	1.8	ug/l	
99-09-2	3-Nitroaniline	ND	4.9	0.86	ug/l	
100-01-6	4-Nitroaniline	ND	4.9	1.1	ug/l	
98-95-3	Nitrobenzene	ND	4.9	0.91	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	4.9	0.66	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	4.9	0.79	ug/l	
85-01-8	Phenanthrene	ND	4.9	0.85	ug/l	
129-00-0	Pyrene	ND	4.9	0.67	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	4.9	0.49	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	24%		14-67%
4165-62-2	Phenol-d5	16%		10-50%
118-79-6	2,4,6-Tribromophenol	71%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	68%		42-108%
321-60-8	2-Fluorobiphenyl	63%		40-106%
1718-51-0	Terphenyl-d14	54%		39-121%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690596.D	1	03/26/17	NG	03/24/17	OP64319	SL3976
Run #2							

	Initial Volume	Final Volume
Run #1	1020 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	9.8	2.5	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	4.9	0.58	ug/l	
95-57-8	2-Chlorophenol	ND	4.9	0.62	ug/l	
120-83-2	2,4-Dichlorophenol	ND	4.9	0.82	ug/l	
105-67-9	2,4-Dimethylphenol	ND	4.9	0.72	ug/l	
51-28-5	2,4-Dinitrophenol	ND	25	4.9	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	9.8	2.0	ug/l	
95-48-7	2-Methylphenol	ND	4.9	0.55	ug/l	
	3&4-Methylphenol	ND	4.9	0.96	ug/l	
88-75-5	2-Nitrophenol	ND	4.9	0.84	ug/l	
100-02-7	4-Nitrophenol	ND	25	4.9	ug/l	
87-86-5	Pentachlorophenol	ND	25	4.9	ug/l	
108-95-2	Phenol	ND	4.9	0.49	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol	ND	4.9	0.95	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	4.9	0.73	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	4.9	0.74	ug/l	
83-32-9	Acenaphthene	ND	4.9	0.61	ug/l	
208-96-8	Acenaphthylene	ND	4.9	0.63	ug/l	
98-86-2	Acetophenone	ND	4.9	0.79	ug/l	
120-12-7	Anthracene	ND	4.9	0.78	ug/l	
1912-24-9	Atrazine	ND	4.9	1.0	ug/l	
100-52-7	Benzaldehyde	ND	25	4.9	ug/l	
56-55-3	Benzo(a)anthracene	ND	4.9	0.75	ug/l	
50-32-8	Benzo(a)pyrene	ND	4.9	0.77	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	4.9	0.76	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	4.9	0.81	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	4.9	0.84	ug/l	
92-52-4	1,1'-Biphenyl	ND	4.9	0.61	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	4.9	0.83	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	4.9	0.98	ug/l	
86-74-8	Carbazole	ND	4.9	0.59	ug/l	
106-47-8	4-Chloroaniline	ND	4.9	0.62	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

COLEGIO DE QUIMICOS DE PUERTO RICO
Rafael Infante Méndez
IQ # 1888
QUIMICO LICENCIADO

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	4.9	0.79	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	4.9	0.72	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	4.9	0.74	ug/l	
91-58-7	2-Chloronaphthalene	ND	4.9	0.49	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	4.9	0.53	ug/l	
218-01-9	Chrysene	ND	4.9	0.83	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	4.9	0.79	ug/l	
132-64-9	Dibenzofuran	ND	4.9	0.59	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	4.9	0.63	ug/l	
84-66-2	Diethyl Phthalate	ND	4.9	0.98	ug/l	
131-11-3	Dimethyl Phthalate	ND	4.9	0.98	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	4.9	0.98	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	4.9	0.98	ug/l	
121-14-2	2,4-Dinitrotoluene	ND	4.9	0.80	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	4.9	0.70	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	4.9	0.98	ug/l	
206-44-0	Fluoranthene	ND	4.9	0.54	ug/l	
86-73-7	Fluorene	ND	4.9	0.69	ug/l	
118-74-1	Hexachlorobenzene	ND	4.9	0.68	ug/l	
87-68-3	Hexachlorobutadiene	ND	4.9	0.49	ug/l	
77-47-4	Hexachlorocyclopentadiene	ND	4.9	1.8	ug/l	
67-72-1	Hexachloroethane	ND	4.9	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	4.9	0.70	ug/l	
78-59-1	Isophorone	ND	4.9	0.76	ug/l	
91-57-6	2-Methylnaphthalene	ND	4.9	0.59	ug/l	
88-74-4	2-Nitroaniline	ND	4.9	1.8	ug/l	
99-09-2	3-Nitroaniline	ND	4.9	0.86	ug/l	
100-01-6	4-Nitroaniline	ND	4.9	1.1	ug/l	
98-95-3	Nitrobenzene	ND	4.9	0.91	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	4.9	0.66	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	4.9	0.79	ug/l	
85-01-8	Phenanthrene	ND	4.9	0.85	ug/l	
129-00-0	Pyrene	ND	4.9	0.67	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	4.9	0.49	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	22%		14-67%
4165-62-2	Phenol-d5	14%		10-50%
118-79-6	2,4,6-Tribromophenol	72%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	65%		42-108%
321-60-8	2-Fluorobiphenyl	59%		40-106%
1718-51-0	Terphenyl-d14	42%		39-121%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690597.D	1	03/26/17	NG	03/24/17	OP64319	SL3976
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	10	2.5	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	5.0	0.59	ug/l	
95-57-8	2-Chlorophenol	ND	5.0	0.63	ug/l	
120-83-2	2,4-Dichlorophenol	ND	5.0	0.84	ug/l	
105-67-9	2,4-Dimethylphenol	ND	5.0	0.74	ug/l	
51-28-5	2,4-Dinitrophenol	ND	25	5.0	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	10	2.0	ug/l	
95-48-7	2-Methylphenol	ND	5.0	0.56	ug/l	
	3&4-Methylphenol	ND	5.0	0.98	ug/l	
88-75-5	2-Nitrophenol	ND	5.0	0.85	ug/l	
100-02-7	4-Nitrophenol	ND	25	5.0	ug/l	
87-86-5	Pentachlorophenol	ND	25	5.0	ug/l	
108-95-2	Phenol	ND	5.0	0.50	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.97	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.74	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.75	ug/l	
83-32-9	Acenaphthene	ND	5.0	0.63	ug/l	
208-96-8	Acenaphthylene	ND	5.0	0.64	ug/l	
98-86-2	Acetophenone	ND	5.0	0.81	ug/l	
120-12-7	Anthracene	ND	5.0	0.80	ug/l	
1912-24-9	Atrazine	ND	5.0	1.0	ug/l	
100-52-7	Benzaldehyde	ND	25	5.0	ug/l	
56-55-3	Benzo(a)anthracene	ND	5.0	0.76	ug/l	
50-32-8	Benzo(a)pyrene	ND	5.0	0.78	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	5.0	0.78	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	5.0	0.82	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	5.0	0.86	ug/l	
92-52-4	1,1'-Biphenyl	ND	5.0	0.62	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	5.0	0.85	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	5.0	1.0	ug/l	
86-74-8	Carbazole	ND	5.0	0.60	ug/l	
106-47-8	4-Chloroaniline	ND	5.0	0.63	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	5.0	0.81	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	5.0	0.73	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	5.0	0.76	ug/l	
91-58-7	2-Chloronaphthalene	ND	5.0	0.50	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	5.0	0.54	ug/l	
218-01-9	Chrysene	ND	5.0	0.85	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	5.0	0.80	ug/l	
132-64-9	Dibenzofuran	ND	5.0	0.60	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	5.0	0.64	ug/l	
84-66-2	Diethyl Phthalate	ND	5.0	1.0	ug/l	
131-11-3	Dimethyl Phthalate	ND	5.0	1.0	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	5.0	1.0	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	5.0	1.0	ug/l	
121-14-2	2,4-Dinitrotoluene	ND	5.0	0.81	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	5.0	0.71	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	5.0	1.0	ug/l	
206-44-0	Fluoranthene	ND	5.0	0.55	ug/l	
86-73-7	Fluorene	ND	5.0	0.70	ug/l	
118-74-1	Hexachlorobenzene	ND	5.0	0.69	ug/l	
87-68-3	Hexachlorobutadiene	ND	5.0	0.50	ug/l	
77-47-4	Hexachlorocyclopentadiene	ND	5.0	1.8	ug/l	
67-72-1	Hexachloroethane	ND	5.0	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	5.0	0.71	ug/l	
78-59-1	Isophorone	ND	5.0	0.78	ug/l	
91-57-6	2-Methylnaphthalene	ND	5.0	0.60	ug/l	
88-74-4	2-Nitroaniline	ND	5.0	1.8	ug/l	
99-09-2	3-Nitroaniline	ND	5.0	0.88	ug/l	
100-01-6	4-Nitroaniline	ND	5.0	1.2	ug/l	
98-95-3	Nitrobenzene	ND	5.0	0.93	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	5.0	0.67	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	5.0	0.81	ug/l	
85-01-8	Phenanthrene	ND	5.0	0.86	ug/l	
129-00-0	Pyrene	ND	5.0	0.68	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	5.0	0.50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	23%		14-67%
4165-62-2	Phenol-d5	17%		10-50%
118-79-6	2,4,6-Tribromophenol	86%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	74%		42-108%
321-60-8	2-Fluorobiphenyl	69%		40-106%
1718-51-0	Terphenyl-d14	44%		39-121%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690600.D	1	03/26/17	NG	03/24/17	OP64319	SL3976
Run #2							

	Initial Volume	Final Volume
Run #1	1040 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	9.6	2.4	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	4.8	0.57	ug/l	
95-57-8	2-Chlorophenol	ND	4.8	0.61	ug/l	
120-83-2	2,4-Dichlorophenol	ND	4.8	0.80	ug/l	
105-67-9	2,4-Dimethylphenol	ND	4.8	0.71	ug/l	
51-28-5	2,4-Dinitrophenol	ND	24	4.8	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	9.6	1.9	ug/l	
95-48-7	2-Methylphenol	ND	4.8	0.54	ug/l	
	3&4-Methylphenol	ND	4.8	0.94	ug/l	
88-75-5	2-Nitrophenol	ND	4.8	0.82	ug/l	
100-02-7	4-Nitrophenol	ND	24	4.8	ug/l	
87-86-5	Pentachlorophenol	ND	24	4.8	ug/l	
108-95-2	Phenol	ND	4.8	0.48	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol	ND	4.8	0.93	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	4.8	0.71	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	4.8	0.72	ug/l	
83-32-9	Acenaphthene	ND	4.8	0.60	ug/l	
208-96-8	Acenaphthylene	ND	4.8	0.61	ug/l	
98-86-2	Acetophenone	ND	4.8	0.78	ug/l	
120-12-7	Anthracene	ND	4.8	0.77	ug/l	
1912-24-9	Atrazine	ND	4.8	0.98	ug/l	
100-52-7	Benzaldehyde	ND	24	4.8	ug/l	
56-55-3	Benzo(a)anthracene	ND	4.8	0.73	ug/l	
50-32-8	Benzo(a)pyrene	ND	4.8	0.75	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	4.8	0.75	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	4.8	0.79	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	4.8	0.82	ug/l	
92-52-4	1,1'-Biphenyl	ND	4.8	0.60	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	4.8	0.81	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	4.8	0.96	ug/l	
86-74-8	Carbazole	ND	4.8	0.58	ug/l	
106-47-8	4-Chloroaniline	ND	4.8	0.61	ug/l	

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	4.8	0.78	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	4.8	0.70	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	4.8	0.73	ug/l	
91-58-7	2-Chloronaphthalene	ND	4.8	0.48	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	4.8	0.52	ug/l	
218-01-9	Chrysene	ND	4.8	0.82	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	4.8	0.77	ug/l	
132-64-9	Dibenzofuran	ND	4.8	0.58	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	4.8	0.62	ug/l	
84-66-2	Diethyl Phthalate	ND	4.8	0.96	ug/l	
131-11-3	Dimethyl Phthalate	ND	4.8	0.96	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	4.8	0.96	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	4.8	0.96	ug/l	
121-14-2	2,4-Dinitrotoluene	ND	4.8	0.78	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	4.8	0.69	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	4.8	0.96	ug/l	
206-44-0	Fluoranthene	ND	4.8	0.53	ug/l	
86-73-7	Fluorene	ND	4.8	0.67	ug/l	
118-74-1	Hexachlorobenzene	ND	4.8	0.67	ug/l	
87-68-3	Hexachlorobutadiene	ND	4.8	0.48	ug/l	
77-47-4	Hexachlorocyclopentadiene	ND	4.8	1.7	ug/l	
67-72-1	Hexachloroethane	ND	4.8	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	4.8	0.69	ug/l	
78-59-1	Isophorone	ND	4.8	0.75	ug/l	
91-57-6	2-Methylnaphthalene	ND	4.8	0.58	ug/l	
88-74-4	2-Nitroaniline	ND	4.8	1.7	ug/l	
99-09-2	3-Nitroaniline	ND	4.8	0.85	ug/l	
100-01-6	4-Nitroaniline	ND	4.8	1.1	ug/l	
98-95-3	Nitrobenzene	ND	4.8	0.90	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	4.8	0.64	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	4.8	0.78	ug/l	
85-01-8	Phenanthrene	ND	4.8	0.83	ug/l	
129-00-0	Pyrene	ND	4.8	0.66	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	4.8	0.48	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	19%		14-67%
4165-62-2	Phenol-d5	12%		10-50%
118-79-6	2,4,6-Tribromophenol	77%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	66%		42-108%
321-60-8	2-Fluorobiphenyl	57%		40-106%
1718-51-0	Terphenyl-d14	39%		39-121%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	Date Sampled:	03/17/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	L0690601.D	1	03/26/17	NG	03/24/17	OP64319	SL3976
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
105-60-2	Caprolactam	ND	10	2.5	ug/l	
59-50-7	4-Chloro-3-methyl Phenol	ND	5.0	0.59	ug/l	
95-57-8	2-Chlorophenol	ND	5.0	0.63	ug/l	
120-83-2	2,4-Dichlorophenol	ND	5.0	0.84	ug/l	
105-67-9	2,4-Dimethylphenol	ND	5.0	0.74	ug/l	
51-28-5	2,4-Dinitrophenol	ND	25	5.0	ug/l	
534-52-1	4,6-Dinitro-o-cresol	ND	10	2.0	ug/l	
95-48-7	2-Methylphenol	ND	5.0	0.56	ug/l	
	3&4-Methylphenol	ND	5.0	0.98	ug/l	
88-75-5	2-Nitrophenol	ND	5.0	0.85	ug/l	
100-02-7	4-Nitrophenol	ND	25	5.0	ug/l	
87-86-5	Pentachlorophenol	ND	25	5.0	ug/l	
108-95-2	Phenol	ND	5.0	0.50	ug/l	
58-90-2	2,3,4,6-Tetrachlorophenol	ND	5.0	0.97	ug/l	
95-95-4	2,4,5-Trichlorophenol	ND	5.0	0.74	ug/l	
88-06-2	2,4,6-Trichlorophenol	ND	5.0	0.75	ug/l	
83-32-9	Acenaphthene	ND	5.0	0.63	ug/l	
208-96-8	Acenaphthylene	ND	5.0	0.64	ug/l	
98-86-2	Acetophenone	ND	5.0	0.81	ug/l	
120-12-7	Anthracene	ND	5.0	0.80	ug/l	
1912-24-9	Atrazine	ND	5.0	1.0	ug/l	
100-52-7	Benzaldehyde	ND	25	5.0	ug/l	
56-55-3	Benzo(a)anthracene	ND	5.0	0.76	ug/l	
50-32-8	Benzo(a)pyrene	ND	5.0	0.78	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	5.0	0.78	ug/l	
191-24-2	Benzo(g,h,i)perylene	ND	5.0	0.82	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	5.0	0.86	ug/l	
92-52-4	1,1'-Biphenyl	ND	5.0	0.62	ug/l	
101-55-3	4-Bromophenyl Phenyl Ether	ND	5.0	0.85	ug/l	
85-68-7	Butyl Benzyl Phthalate	ND	5.0	1.0	ug/l	
86-74-8	Carbazole	ND	5.0	0.60	ug/l	
106-47-8	4-Chloroaniline	ND	5.0	0.63	ug/l	

ND = Not detected MDL = Method Detection Limit
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J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	**Date Sampled:**	03/17/17
Matrix:	AQ - Field Blank Water	**Date Received:**	03/18/17
Method:	SW846 8270D SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
111-91-1	bis(2-Chloroethoxy)methane	ND	5.0	0.81	ug/l	
111-44-4	bis(2-Chloroethyl)ether	ND	5.0	0.73	ug/l	
108-60-1	bis(2-Chloroisopropyl)ether	ND	5.0	0.76	ug/l	
91-58-7	2-Chloronaphthalene	ND	5.0	0.50	ug/l	
7005-72-3	4-Chlorophenyl Phenyl Ether	ND	5.0	0.54	ug/l	
218-01-9	Chrysene	ND	5.0	0.85	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	5.0	0.80	ug/l	
132-64-9	Dibenzofuran	ND	5.0	0.60	ug/l	
91-94-1	3,3'-Dichlorobenzidine	ND	5.0	0.64	ug/l	
84-66-2	Diethyl Phthalate	ND	5.0	1.0	ug/l	
131-11-3	Dimethyl Phthalate	ND	5.0	1.0	ug/l	
84-74-2	Di-n-butyl Phthalate	ND	5.0	1.0	ug/l	
117-84-0	Di-n-octyl Phthalate	ND	5.0	1.0	ug/l	
121-14-2	2,4-Dinitrotoluene	ND	5.0	0.81	ug/l	
606-20-2	2,6-Dinitrotoluene	ND	5.0	0.71	ug/l	
117-81-7	bis(2-Ethylhexyl)phthalate	ND	5.0	1.0	ug/l	
206-44-0	Fluoranthene	ND	5.0	0.55	ug/l	
86-73-7	Fluorene	ND	5.0	0.70	ug/l	
118-74-1	Hexachlorobenzene	ND	5.0	0.69	ug/l	
87-68-3	Hexachlorobutadiene	ND	5.0	0.50	ug/l	
77-47-4	Hexachlorocyclopentadiene	ND	5.0	1.8	ug/l	
67-72-1	Hexachloroethane	ND	5.0	1.6	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	5.0	0.71	ug/l	
78-59-1	Isophorone	ND	5.0	0.78	ug/l	
91-57-6	2-Methylnaphthalene	ND	5.0	0.60	ug/l	
88-74-4	2-Nitroaniline	ND	5.0	1.8	ug/l	
99-09-2	3-Nitroaniline	ND	5.0	0.88	ug/l	
100-01-6	4-Nitroaniline	ND	5.0	1.2	ug/l	
98-95-3	Nitrobenzene	ND	5.0	0.93	ug/l	
621-64-7	N-Nitrosodi-n-propylamine	ND	5.0	0.67	ug/l	
86-30-6	N-Nitrosodiphenylamine	ND	5.0	0.81	ug/l	
85-01-8	Phenanthrene	ND	5.0	0.86	ug/l	
129-00-0	Pyrene	ND	5.0	0.68	ug/l	
95-94-3	1,2,4,5-Tetrachlorobenzene	ND	5.0	0.50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
367-12-4	2-Fluorophenol	29%		14-67%
4165-62-2	Phenol-d5	18%		10-50%
118-79-6	2,4,6-Tribromophenol	79%		33-118%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	Date Sampled:	03/17/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846 8270D SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

ABN TCL Special List

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	77%		42-108%
321-60-8	2-Fluorobiphenyl	72%		40-106%
1718-51-0	Terphenyl-d14	53%		39-121%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

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Report of Analysis

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	Date Sampled:	03/16/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098461.D	1	03/31/17	FS	03/23/17	OP64303	SW4371
Run #2	U060594.D	1	03/24/17	NJ	03/23/17	OP64303	SU2655

	Initial Volume	Final Volume
Run #1	1020 ml	1.0 ml
Run #2	1020 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.20	0.039	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.20	0.039	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.20	0.039	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.20	0.039	ug/l	
218-01-9	Chrysene	ND	0.20	0.039	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.20	0.039	ug/l	
123-91-1	1,4-Dioxane	ND [a]	0.29	0.15	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.20	0.039	ug/l	
90-12-0	1-Methylnaphthalene	ND	0.98	0.39	ug/l	
91-57-6	2-Methylnaphthalene	ND	0.98	0.39	ug/l	
91-20-3	Naphthalene	ND	0.98	0.39	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	54% [b]	61% [b]	42-108%
321-60-8	2-Fluorobiphenyl	51% [b]	54% [b]	40-106%
1718-51-0	Terphenyl-d14	45% [b]	64% [b]	39-121%

(a) Result is from Run# 2
(b) Surrogate recoveries corrected for actual spike amount.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098462.D	1	03/31/17	FS	03/23/17	OP64303	SW4371
Run #2	U060595.D	1	03/24/17	NJ	03/23/17	OP64303	SU2655

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2	1000 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.20	0.040	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.20	0.040	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.20	0.040	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.20	0.040	ug/l	
218-01-9	Chrysene	ND	0.20	0.040	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.20	0.040	ug/l	
123-91-1	1,4-Dioxane	15.9 [a]	0.30	0.15	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.20	0.040	ug/l	
90-12-0	1-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-57-6	2-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-20-3	Naphthalene	ND	1.0	0.40	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	61% [b]	73% [b]	42-108%
321-60-8	2-Fluorobiphenyl	65% [b]	60% [b]	40-106%
1718-51-0	Terphenyl-d14	50% [b]	69% [b]	39-121%

(a) Result is from Run# 2
(b) Surrogate recoveries corrected for actual spike amount.

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098463.D	1	03/31/17	FS	03/23/17	OP64303	SW4371
Run #2	U060596.D	1	03/24/17	NJ	03/23/17	OP64303	SU2655

	Initial Volume	Final Volume
Run #1	1020 ml	1.0 ml
Run #2	1020 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.20	0.039	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.20	0.039	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.20	0.039	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.20	0.039	ug/l	
218-01-9	Chrysene	ND	0.20	0.039	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.20	0.039	ug/l	
123-91-1	1,4-Dioxane	18.4 [a]	0.29	0.15	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.20	0.039	ug/l	
90-12-0	1-Methylnaphthalene	ND	0.98	0.39	ug/l	
91-57-6	2-Methylnaphthalene	ND	0.98	0.39	ug/l	
91-20-3	Naphthalene	ND	0.98	0.39	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	65% [b]	71% [b]	42-108%
321-60-8	2-Fluorobiphenyl	68% [b]	60% [b]	40-106%
1718-51-0	Terphenyl-d14	50% [b]	67% [b]	39-121%

(a) Result is from Run# 2
(b) Surrogate recoveries corrected for actual spike amount.

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098464.D	1	03/31/17	FS	03/23/17	OP64303	SW4371
Run #2	U060597.D	1	03/24/17	NJ	03/23/17	OP64303	SU2655

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2	1000 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.20	0.040	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.20	0.040	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.20	0.040	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.20	0.040	ug/l	
218-01-9	Chrysene	ND	0.20	0.040	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.20	0.040	ug/l	
123-91-1	1,4-Dioxane	3.0 [a]	0.30	0.15	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.20	0.040	ug/l	
90-12-0	1-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-57-6	2-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-20-3	Naphthalene	0.40	1.0	0.40	ug/l	J

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	60% [b]	70% [b]	42-108%
321-60-8	2-Fluorobiphenyl	66% [b]	63% [b]	40-106%
1718-51-0	Terphenyl-d14	51% [b]	70% [b]	39-121%

(a) Result is from Run# 2
(b) Surrogate recoveries corrected for actual spike amount.

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098465.D	1	03/31/17	FS	03/23/17	OP64303	SW4371
Run #2	U060598.D	1	03/24/17	NJ	03/23/17	OP64303	SU2655

	Initial Volume	Final Volume
Run #1	1040 ml	1.0 ml
Run #2	1040 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.19	0.038	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.19	0.038	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.19	0.038	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.19	0.038	ug/l	
218-01-9	Chrysene	ND	0.19	0.038	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.19	0.038	ug/l	
123-91-1	1,4-Dioxane	2.6 [a]	0.29	0.14	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.19	0.038	ug/l	
90-12-0	1-Methylnaphthalene	ND	0.96	0.38	ug/l	
91-57-6	2-Methylnaphthalene	ND	0.96	0.38	ug/l	
91-20-3	Naphthalene	ND	0.96	0.38	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	60% [b]	70% [b]	42-108%
321-60-8	2-Fluorobiphenyl	67% [b]	64% [b]	40-106%
1718-51-0	Terphenyl-d14	51% [b]	74% [b]	39-121%

(a) Result is from Run# 2
(b) Surrogate recoveries corrected for actual spike amount.

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ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098350.D	1	03/28/17	FS	03/24/17	OP64320	SW4368
Run #2	U060624.D	1	03/25/17	NJ	03/24/17	OP64320	SU2656

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2	1000 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.20	0.040	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.20	0.040	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.20	0.040	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.20	0.040	ug/l	
218-01-9	Chrysene	ND	0.20	0.040	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.20	0.040	ug/l	
123-91-1	1,4-Dioxane	1.2 [a]	0.30	0.15	ug/l	B
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.20	0.040	ug/l	
90-12-0	1-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-57-6	2-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-20-3	Naphthalene	ND	1.0	0.40	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	53%	61% [b]	42-108%
321-60-8	2-Fluorobiphenyl	60%	56% [b]	40-106%
1718-51-0	Terphenyl-d14	42%	73% [b]	39-121%

(a) Result is from Run# 2
(b) Surrogate recoveries corrected for actual spike amount.

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Rafael Infante Méndez
IQ # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098351.D	1	03/28/17	FS	03/24/17	OP64320	SW4368
Run #2	U060625.D	1	03/25/17	NJ	03/24/17	OP64320	SU2656

	Initial Volume	Final Volume
Run #1	1020 ml	1.0 ml
Run #2	1020 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.20	0.039	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.20	0.039	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.20	0.039	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.20	0.039	ug/l	
218-01-9	Chrysene	ND	0.20	0.039	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.20	0.039	ug/l	
123-91-1	1,4-Dioxane	1.5 [a]	0.29	0.15	ug/l	B
123-91-1	1,4-Dioxane	ND	0.29	0.15	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.20	0.039	ug/l	
90-12-0	1-Methylnaphthalene	ND	0.98	0.39	ug/l	
91-57-6	2-Methylnaphthalene	ND	0.98	0.39	ug/l	
91-20-3	Naphthalene	ND	0.98	0.39	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	58%	66% [b]	42-108%
321-60-8	2-Fluorobiphenyl	56%	58% [b]	40-106%
1718-51-0	Terphenyl-d14	43%	64% [b]	39-121%

(a) Result is from Run# 2
(b) Surrogate recoveries corrected for actual spike amount.

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098352.D	1	03/28/17	FS	03/24/17	OP64320	SW4368
Run #2	U060657.D	10	03/29/17	NJ	03/24/17	OP64320	SU2657
Run #3 [a]	W098509.D	1	04/03/17	FS	03/31/17	OP64441	SW4372

	Initial Volume	Final Volume
Run #1	1020 ml	1.0 ml
Run #2	1020 ml	1.0 ml
Run #3	1040 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.20	0.039	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.20	0.039	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.20	0.039	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.20	0.039	ug/l	
218-01-9	Chrysene	ND	0.20	0.039	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.20	0.039	ug/l	
123-91-1	1,4-Dioxane	ND	0.29	0.15	ug/l	
123-91-1	1,4-Dioxane	103 [b]	2.9	1.5	ug/l	B
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.20	0.039	ug/l	
90-12-0	1-Methylnaphthalene	ND	0.98	0.39	ug/l	
91-57-6	2-Methylnaphthalene	ND	0.98	0.39	ug/l	
91-20-3	Naphthalene	ND	0.98	0.39	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Run# 3	Limits
4165-60-0	Nitrobenzene-d5	52%	55% [c]	74%	42-108%
321-60-8	2-Fluorobiphenyl	51%	51% [c]	80%	40-106%
1718-51-0	Terphenyl-d14	31% [d]	46% [c]	77%	39-121%

(a) Confirmation run.
(b) Result is from Run# 2
(c) Surrogate recoveries corrected for actual spike amount.
(d) Outside control limits due to matrix interference. Confirmed by re-extraction and reanalysis beyond hold time.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098353.D	1	03/28/17	FS	03/24/17	OP64320	SW4368
Run #2	U060658.D	25	03/29/17	NJ	03/24/17	OP64320	SU2657
Run #3 [a]	W098510.D	1	04/03/17	FS	03/31/17	OP64441	SW4372

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2	1000 ml	1.0 ml
Run #3	1020 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.20	0.040	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.20	0.040	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.20	0.040	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.20	0.040	ug/l	
218-01-9	Chrysene	ND	0.20	0.040	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.20	0.040	ug/l	
123-91-1	1,4-Dioxane	ND	0.30	0.15	ug/l	
123-91-1	1,4-Dioxane	313 [b]	7.5	3.8	ug/l	B
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.20	0.040	ug/l	
90-12-0	1-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-57-6	2-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-20-3	Naphthalene	ND	1.0	0.40	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Run# 3	Limits
4165-60-0	Nitrobenzene-d5	57%	0% [c]	74%	42-108%
321-60-8	2-Fluorobiphenyl	63%	0% [c]	83%	40-106%
1718-51-0	Terphenyl-d14	35% [d]	0% [c]	67%	39-121%

(a) Confirmation run.
(b) Result is from Run# 2
(c) Outside control limits due to dilution.
(d) Outside control limits due to matrix interference. Confirmed by re-extraction and reanalysis beyond hold time.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098354.D	1	03/28/17	FS	03/24/17	OP64320	SW4368
Run #2	U060628.D	1	03/25/17	NJ	03/24/17	OP64320	SU2656
Run #3 [a]	W098511.D	1	04/03/17	FS	03/31/17	OP64441	SW4372

	Initial Volume	Final Volume
Run #1	1040 ml	1.0 ml
Run #2	1040 ml	1.0 ml
Run #3	1040 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.19	0.038	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.19	0.038	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.19	0.038	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.19	0.038	ug/l	
218-01-9	Chrysene	ND	0.19	0.038	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.19	0.038	ug/l	
123-91-1	1,4-Dioxane	15.8 [b]	0.29	0.14	ug/l	B
123-91-1	1,4-Dioxane	ND	0.29	0.14	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.19	0.038	ug/l	
90-12-0	1-Methylnaphthalene	ND	0.96	0.38	ug/l	
91-57-6	2-Methylnaphthalene	ND	0.96	0.38	ug/l	
91-20-3	Naphthalene	ND	0.96	0.38	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Run# 3	Limits
4165-60-0	Nitrobenzene-d5	49%	59% [c]	71%	42-108%
321-60-8	2-Fluorobiphenyl	48%	51% [c]	77%	40-106%
1718-51-0	Terphenyl-d14	29% [d]	40% [c]	71%	39-121%

(a) Confirmation run.
(b) Result is from Run# 2
(c) Surrogate recoveries corrected for actual spike amount.
(d) Outside control limits due to matrix interference. Confirmed by re-extraction and reanalysis beyond hold time.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	**Date Sampled:**	03/17/17
Matrix:	AQ - Field Blank Water	**Date Received:**	03/18/17
Method:	SW846 8270D BY SIM SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	W098355.D	1	03/28/17	FS	03/24/17	OP64320	SW4368
Run #2	U060629.D	1	03/25/17	NJ	03/24/17	OP64320	SU2656

	Initial Volume	Final Volume
Run #1	1000 ml	1.0 ml
Run #2	1000 ml	1.0 ml

BN Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
56-55-3	Benzo(a)anthracene	ND	0.20	0.040	ug/l	
50-32-8	Benzo(a)pyrene	ND	0.20	0.040	ug/l	
205-99-2	Benzo(b)fluoranthene	ND	0.20	0.040	ug/l	
207-08-9	Benzo(k)fluoranthene	ND	0.20	0.040	ug/l	
218-01-9	Chrysene	ND	0.20	0.040	ug/l	
53-70-3	Dibenzo(a,h)anthracene	ND	0.20	0.040	ug/l	
123-91-1	1,4-Dioxane	ND [a]	0.30	0.15	ug/l	
123-91-1	1,4-Dioxane	ND	0.30	0.15	ug/l	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.20	0.040	ug/l	
90-12-0	1-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-57-6	2-Methylnaphthalene	ND	1.0	0.40	ug/l	
91-20-3	Naphthalene	ND	1.0	0.40	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	63%	74% [b]	42-108%
321-60-8	2-Fluorobiphenyl	66%	65% [b]	40-106%
1718-51-0	Terphenyl-d14	38%	56% [b]	39-121%

(a) Result is from Run# 2
(b) Surrogate recoveries corrected for actual spike amount.

LIBRE ASOCIADO DE PUERTO RICO
dael Infante
Méndez
IC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	Date Sampled:	03/16/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018135.D	1	04/12/17	MG	03/23/17	OP64317	GNN913
Run #2							

	Initial Volume	Final Volume
Run #1	1050 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	99.1	190	76	ug/l	J
	C9-C18 Aliphatics	ND	190	48	ug/l	
	C19-C36 Aliphatics	ND	190	48	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	70%		40-140%
580-13-2	2-Bromonaphthalene	88%		40-140%
84-15-1	o-Terphenyl	67%		40-140%
321-60-8	2-Fluorobiphenyl	84%		40-140%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018136.D	1	04/12/17	MG	03/23/17	OP64317	GNN913
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	152	200	80	ug/l	J
	C9-C18 Aliphatics	ND	200	50	ug/l	
	C19-C36 Aliphatics	ND	200	50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	69%		40-140%
580-13-2	2-Bromonaphthalene	94%		40-140%
84-15-1	o-Terphenyl	67%		40-140%
321-60-8	2-Fluorobiphenyl	89%		40-140%

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1868

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018137.D	1	04/12/17	MG	03/23/17	OP64317	GNN913
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	147	200	80	ug/l	J
	C9-C18 Aliphatics	ND	200	50	ug/l	
	C19-C36 Aliphatics	ND	200	50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	60%		40-140%
580-13-2	2-Bromonaphthalene	78%		40-140%
84-15-1	o-Terphenyl	54%		40-140%
321-60-8	2-Fluorobiphenyl	74%		40-140%

LIBRE ASOCIADO DE PUERTO RICO
...fael Infante
Méndez
IC # 1888
...UIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018138.D	1	04/12/17	MG	03/23/17	OP64317	GNN913
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	165	200	80	ug/l	J
	C9-C18 Aliphatics	95.3	200	50	ug/l	J
	C19-C36 Aliphatics	50.2	200	50	ug/l	J

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	42%		40-140%
580-13-2	2-Bromonaphthalene	78%		40-140%
84-15-1	o-Terphenyl	40%		40-140%
321-60-8	2-Fluorobiphenyl	74%		40-140%

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IQ # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018139.D	1	04/12/17	MG	03/23/17	OP64317	GNN913
Run #2							

	Initial Volume	Final Volume
Run #1	1050 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	133	190	76	ug/l	J
	C9-C18 Aliphatics	ND	190	48	ug/l	
	C19-C36 Aliphatics	ND	190	48	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	45%		40-140%
580-13-2	2-Bromonaphthalene	73%		40-140%
84-15-1	o-Terphenyl	42%		40-140%
321-60-8	2-Fluorobiphenyl	70%		40-140%

ESTADO LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1086
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018141.D	1	04/12/17	MG	03/23/17	OP64317	GNN913
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	152	200	80	ug/l	J
	C9-C18 Aliphatics	ND	200	50	ug/l	
	C19-C36 Aliphatics	ND	200	50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	55%		40-140%
580-13-2	2-Bromonaphthalene	95%		40-140%
84-15-1	o-Terphenyl	66%		40-140%
321-60-8	2-Fluorobiphenyl	88%		40-140%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018207.D	1	04/15/17	MG	03/23/17	OP64317	GNN916
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	ND	200	80	ug/l	
	C9-C18 Aliphatics	ND	200	50	ug/l	
	C19-C36 Aliphatics	ND	200	50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	42%		40-140%
580-13-2	2-Bromonaphthalene	78%		40-140%
84-15-1	o-Terphenyl	64%		40-140%
321-60-8	2-Fluorobiphenyl	75%		40-140%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018230.D	1	04/17/17	MG	03/23/17	OP64317	GNN917
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	ND	200	80	ug/l	
	C9-C18 Aliphatics	ND	200	50	ug/l	
	C19-C36 Aliphatics	ND	200	50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	40%		40-140%
580-13-2	2-Bromonaphthalene	84%		40-140%
84-15-1	o-Terphenyl	65%		40-140%
321-60-8	2-Fluorobiphenyl	79%		40-140%

LIBRE ASOCIADO DE PUERTO RICO
Manuel Infante Méndez
1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018179.D	1	04/13/17	MG	03/23/17	OP64317	GNN914
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	149	200	80	ug/l	J
	C9-C18 Aliphatics	ND	200	50	ug/l	
	C19-C36 Aliphatics	ND	200	50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	57%		40-140%
580-13-2	2-Bromonaphthalene	91%		40-140%
84-15-1	o-Terphenyl	73%		40-140%
321-60-8	2-Fluorobiphenyl	86%		40-140%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018180.D	1	04/13/17	MG	03/23/17	OP64317	GNN914
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	134	200	80	ug/l	J
	C9-C18 Aliphatics	ND	200	50	ug/l	
	C19-C36 Aliphatics	ND	200	50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	60%		40-140%
580-13-2	2-Bromonaphthalene	89%		40-140%
84-15-1	o-Terphenyl	66%		40-140%
321-60-8	2-Fluorobiphenyl	85%		40-140%

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	Date Sampled:	03/17/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	MADEP EPH REV 1.1 SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	NN018209.D	1	04/15/17	MG	03/23/17	OP64317	GNN916
Run #2							

	Initial Volume	Final Volume
Run #1	1000 ml	2.0 ml
Run #2		

MAEPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C11-C22 Aromatics (Unadj.)	ND	200	80	ug/l	
	C9-C18 Aliphatics	ND	200	50	ug/l	
	C19-C36 Aliphatics	ND	200	50	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
3386-33-2	1-Chlorooctadecane	43%		40-140%
580-13-2	2-Bromonaphthalene	71%		40-140%
84-15-1	o-Terphenyl	57%		40-140%
321-60-8	2-Fluorobiphenyl	68%		40-140%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

4.1
4

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	Date Sampled:	03/16/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	TT381599.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [a]	ND	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane	ND	0.040	0.0088	ug/l	
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [a]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [a]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde	ND	0.080	0.011	ug/l	
53494-70-5	Endrin ketone [a]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	111%		42-127%
2051-24-3	Decachlorobiphenyl	121%		27-127%

(a) Associated BS recovery outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	TT381600.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [a]	ND	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane [b]	0.013	0.040	0.0088	ug/l	J
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [a]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [a]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde	ND	0.080	0.011	ug/l	
53494-70-5	Endrin ketone [a]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	106%		42-127%
2051-24-3	Decachlorobiphenyl	130%		27-127%

(a) Associated BS recovery outside control limits.
(b) Primary and confirmation results differ by more than 40%. Lower value reported due to possible coelution.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	TT381601.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [b]	ND	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane	0.017	0.040	0.0088	ug/l	J
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [b]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [b]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde	ND	0.080	0.011	ug/l	
53494-70-5	Endrin ketone [b]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	115%		42-127%
2051-24-3	Decachlorobiphenyl	133%		27-127%

(a) All hits confirmed by dual column analysis.
(b) Associated BS recovery outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	KK82632.D	1	04/03/17	MV	03/23/17	OP64318	GKK2644
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	0.025	0.040	0.0095	ug/l	J
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane	ND	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane	ND	0.040	0.0088	ug/l	
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde	ND	0.080	0.011	ug/l	
53494-70-5	Endrin ketone	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor [b]	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	98%		42-127%
2051-24-3	Decachlorobiphenyl	113%		27-127%

(a) All hits confirmed by dual column analysis.
(b) Associated CCV outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	TT381607.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [b]	0.041	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane	ND	0.040	0.0088	ug/l	
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [b]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [b]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde [c]	ND	0.080	0.011	ug/l	
53494-70-5	Endrin ketone [b]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	110%		42-127%
2051-24-3	Decachlorobiphenyl	127%		27-127%

(a) All hits confirmed by dual column analysis.
(b) Associated BS recovery outside control limits.
(c) Associated CCV outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	TT381608.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [a]	ND	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane	ND	0.040	0.0088	ug/l	
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [a]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [a]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde [b]	ND	0.080	0.080	ug/l	
53494-70-5	Endrin ketone [a]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	115%		42-127%
2051-24-3	Decachlorobiphenyl	115%		27-127%

(a) Associated BS recovery outside control limits.
(b) Associated CCV outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	TT381609.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [a]	ND	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane	ND	0.040	0.0088	ug/l	
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [a]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [a]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde [b]	ND	0.080	0.080	ug/l	
53494-70-5	Endrin ketone [a]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	107%		42-127%
2051-24-3	Decachlorobiphenyl	102%		27-127%

(a) Associated BS recovery outside control limits.
(b) Associated CCV outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	TT381610.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [b]	0.046	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane	ND	0.040	0.0088	ug/l	
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [b]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [b]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde [c]	ND	0.080	0.080	ug/l	
53494-70-5	Endrin ketone [b]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	132%		42-127%
2051-24-3	Decachlorobiphenyl	153%		27-127%

(a) All hits confirmed by dual column analysis.
(b) Associated BS recovery outside control limits.
(c) Associated CCV outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	TT381611.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [b]	0.025	0.040	0.0077	ug/l	J
5103-74-2	gamma-Chlordane	ND	0.040	0.0088	ug/l	
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [b]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [b]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde [c]	ND	0.080	0.080	ug/l	
53494-70-5	Endrin ketone [b]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	116%		42-127%
2051-24-3	Decachlorobiphenyl	135% [d]		27-127%

(a) All hits confirmed by dual column analysis.
(b) Associated BS recovery outside control limits.
(c) Associated CCV outside control limits.
(d) Outside control limits.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846 8081B SW846 3510C	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	TT381612.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [a]	ND	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane	ND	0.040	0.0088	ug/l	
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [a]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [a]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde [b]	ND	0.080	0.080	ug/l	
53494-70-5	Endrin ketone [a]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	115%		42-127%
2051-24-3	Decachlorobiphenyl	129% [c]		27-127%

(a) Associated BS recovery outside control limits.
(b) Associated CCV outside control limits.
(c) Outside control limits.

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Méndez
1088
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	**Date Sampled:**	03/17/17
Matrix:	AQ - Field Blank Water	**Date Received:**	03/18/17
Method:	SW846 8081B SW846 3510C	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	TT381613.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
Run #2							

	Initial Volume	Final Volume
Run #1	250 ml	5.0 ml
Run #2		

Pesticide TCL List

CAS No.	Compound	Result	RL	MDL	Units	Q
309-00-2	Aldrin	ND	0.040	0.011	ug/l	
319-84-6	alpha-BHC	ND	0.040	0.0087	ug/l	
319-85-7	beta-BHC	ND	0.040	0.010	ug/l	
319-86-8	delta-BHC	ND	0.040	0.0095	ug/l	
58-89-9	gamma-BHC (Lindane)	ND	0.040	0.0089	ug/l	
5103-71-9	alpha-Chlordane [a]	ND	0.040	0.0077	ug/l	
5103-74-2	gamma-Chlordane	ND	0.040	0.0088	ug/l	
60-57-1	Dieldrin	ND	0.040	0.0095	ug/l	
72-54-8	4,4'-DDD [a]	ND	0.080	0.020	ug/l	
72-55-9	4,4'-DDE	ND	0.080	0.020	ug/l	
50-29-3	4,4'-DDT	ND	0.080	0.020	ug/l	
72-20-8	Endrin	ND	0.080	0.0084	ug/l	
1031-07-8	Endosulfan sulfate [a]	ND	0.080	0.0063	ug/l	
7421-93-4	Endrin aldehyde [b]	ND	0.080	0.011	ug/l	
53494-70-5	Endrin ketone [a]	ND	0.080	0.0062	ug/l	
959-98-8	Endosulfan-I	ND	0.040	0.0064	ug/l	
33213-65-9	Endosulfan-II	ND	0.040	0.0060	ug/l	
76-44-8	Heptachlor	ND	0.040	0.010	ug/l	
1024-57-3	Heptachlor epoxide	ND	0.040	0.0081	ug/l	
72-43-5	Methoxychlor	ND	0.080	0.020	ug/l	
8001-35-2	Toxaphene	ND	2.0	0.86	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
877-09-8	Tetrachloro-m-xylene	133% [c]		42-127%
2051-24-3	Decachlorobiphenyl	120%		27-127%

(a) Associated BS recovery outside control limits.
(b) Associated CCV outside control limits.
(c) Outside control limits.

ND = Not detected MDL = Method Detection Limit
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E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

Matrix Spike/Matrix Spike Duplicate Summary

Job Number: FA42192
Account: AMANYWP Anderson, Mulholland & Associates
Project: BMSMC, Humacao, PR

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP64318-MS	TT381602.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
OP64318-MSD	TT381603.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932
FA42192-3 [a]	TT381601.D	1	03/28/17	MV	03/23/17	OP64318	GTT1932

The QC reported here applies to the following samples: **Method:** SW846 8081B

FA42192-1, FA42192-2, FA42192-3, FA42192-6, FA42192-7, FA42192-8, FA42192-11, FA42192-12, FA42192-13, FA42192-14, FA42192-15

CAS No.	Compound	FA42192-3 ug/l	Q	Spike ug/l	MS ug/l	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
309-00-2	Aldrin	ND		1	1.2	120	1	1.3	130*	8	61-126/21
319-84-6	alpha-BHC	ND		1	1.2	120	1	1.3	130*	8	66-129/23
319-85-7	beta-BHC	ND		1	1.2	120	1	1.3	130	8	66-132/23
319-86-8	delta-BHC	ND		1	1.3	130	1	1.4	140	7	41-142/27
58-89-9	gamma-BHC (Lindane)	ND		1	1.2	120	1	1.3	130	8	68-132/22
5103-71-9	alpha-Chlordane	ND		1	1.4	140*	1	1.5	150*	7	66-131/24
5103-74-2	gamma-Chlordane	0.017	J	1	1.7	167*	1	2.0	197*	16	68-128/21
60-57-1	Dieldrin	ND		1	1.3	130	1	1.4	140*	7	66-138/22
72-54-8	4,4'-DDD	ND		1	1.4	140*	1	1.5	150*	7	63-138/24
72-55-9	4,4'-DDE	ND		1	1.3	130	1	1.4	140*	7	59-133/23
50-29-3	4,4'-DDT	ND		1	1.4	140	1	1.5	150*	7	55-145/27
72-20-8	Endrin	ND		1	1.4	140	1	1.5	150*	7	71-147/23
1031-07-8	Endosulfan sulfate	ND		1	1.4	140*	1	1.5	150*	7	64-128/26
7421-93-4	Endrin aldehyde	ND		1	1.3	122	1	1.3	122	0	60-130/22
53494-70-5	Endrin ketone	ND		1	1.5	150*	1	1.5	150*	0	66-137/25
959-98-8	Endosulfan-I	ND		1	1.2	120	1	1.2	120	0	66-133/21
33213-65-9	Endosulfan-II	ND		1	1.3	130	1	1.4	140*	7	65-133/22
76-44-8	Heptachlor	ND		1	1.2	120	1	1.3	130	8	63-130/23
1024-57-3	Heptachlor epoxide	ND		1	1.2	120	1	1.3	130*	8	67-129/23
72-43-5	Methoxychlor	ND		1	1.4	140*	1	1.5	150*	7	60-136/25

CAS No.	Surrogate Recoveries	MS	MSD	FA42192-3	Limits
877-09-8	Tetrachloro-m-xylene	107%	117%	115%	42-127%
2051-24-3	Decachlorobiphenyl	135%*	133%*	133%*	27-127%

(a) All hits confirmed by dual column analysis.

12.4.1 12

* = Outside of Control Limits.

SGS Accutest

Report of Analysis

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	Date Sampled:	03/16/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	MADEP VPH REV 1.1	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078895.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	ND	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	95%		70-130%
460-00-4	BFB	95%		70-130%

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Méndez
IC # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP VPH REV 1.1	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078896.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	ND	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	95%		70-130%
460-00-4	BFB	95%		70-130%

LIBRE ASOCIADO DE PUERTO RICO
Manuel Infante Méndez
IQ # 1886
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP VPH REV 1.1	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078897.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	35.6	100	35	ug/l	J

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	96%		70-130%
460-00-4	BFB	97%		70-130%

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1888

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP VPH REV 1.1	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078898.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2	UV078931.D	500	03/27/17	AJC	n/a	n/a	GUV4187

	Purge Volume
Run #1	5.0 ml
Run #2	5.0 ml

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	38.4	100	35	ug/l	J
	C9- C12 Aliphatics (Unadj.)	75400 [a]	50000	18000	ug/l	
	C9- C10 Aromatics (Unadj.)	262	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	124%	95%	70-130%
460-00-4	BFB	98%	98%	70-130%

(a) Result is from Run# 2

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IQ # 1[illegible]
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP VPH REV 1.1	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078899.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	ND	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	94%		70-130%
460-00-4	BFB	92%		70-130%

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IQ # 1888
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	MADEP VPH REV 1.1	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078900.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	ND	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	95%		70-130%
460-00-4	BFB	97%		70-130%

LIBRE ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
IC # 1868
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP VPH REV 1.1	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078901.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	ND	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	93%		70-130%
460-00-4	BFB	95%		70-130%

LIBRE ASOCIADO DE PUERTO RICO
...uel Infante
Méndez
IC # 1[illegible]
QUIMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP VPH REV 1.1	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078904.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	ND	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	94%		70-130%
460-00-4	BFB	94%		70-130%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	MADEP VPH REV 1.1	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078905.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	ND	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	91%		70-130%
460-00-4	BFB	91%		70-130%

ABOGADO DE PUERTO RICO
tael Infante
Méndez
1888
QUÍMICO LICENCIADO

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	MADEP VPH REV 1.1	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078906.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	ND	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	92%		70-130%
460-00-4	BFB	92%		70-130%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	Date Sampled:	03/17/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	MADEP VPH REV 1.1	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	UV078907.D	1	03/26/17	AJC	n/a	n/a	GUV4186
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

MADEP VPH List

CAS No.	Compound	Result	RL	MDL	Units	Q
	C5- C8 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C12 Aliphatics (Unadj.)	ND	100	35	ug/l	
	C9- C10 Aromatics (Unadj.)	ND	100	35	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
460-00-4	BFB	93%		70-130%
460-00-4	BFB	94%		70-130%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	FB031617		
Lab Sample ID:	FA42192-1	Date Sampled:	03/16/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846-8015C (DAI)	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109122.D	1	03/24/17	ANJ	n/a	n/a	N:GGH5693
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	79%		56-145%
111-27-3	Hexanol	78%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-2	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846-8015C (DAI)	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109123.D	1	03/24/17	ANJ	n/a	n/a	N:GGH5693
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	69%		56-145%
111-27-3	Hexanol	68%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-34		
Lab Sample ID:	FA42192-3	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846-8015C (DAI)	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109124.D	1	03/24/17	ANJ	n/a	n/a	N:GGH5693
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	79%		56-145%
111-27-3	Hexanol	73%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

LABORATORIO ASOCIADO DE PUERTO RICO
Rafael Infante Méndez
ID # 1008

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-32		
Lab Sample ID:	FA42192-6	**Date Sampled:**	03/16/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846-8015C (DAI)	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109125.D	1	03/24/17	ANJ	n/a	n/a	N:GGH5693
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	78%		56-145%
111-27-3	Hexanol	75%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

LABORATORIO QUÍMICO ASOCIADO DE PUERTO RICO
Manuel Infante Méndez
IQ # 1582

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-36		
Lab Sample ID:	FA42192-7	Date Sampled:	03/16/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846-8015C (DAI)	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109126.D	1	03/24/17	ANJ	n/a	n/a	N:GGH5693
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	81%		56-145%
111-27-3	Hexanol	76%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

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Rafael Infante Méndez
IC # [illegible]

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1		
Lab Sample ID:	FA42192-8	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846-8015C (DAI)	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109127.D	1	03/24/17	ANJ	n/a	n/a	N:GGH5693
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	80%		56-145%
111-27-3	Hexanol	78%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	UP-1DUP		
Lab Sample ID:	FA42192-11	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846-8015C (DAI)	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109128.D	1	03/24/17	ANJ	n/a	n/a	N:GGH5693
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	77%		56-145%
111-27-3	Hexanol	76%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-28		
Lab Sample ID:	FA42192-12	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846-8015C (DAI)	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109129.D	1	03/24/17	ANJ	n/a	n/a	N:GGH5693
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	75%		56-145%
111-27-3	Hexanol	75%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	UP-2		
Lab Sample ID:	FA42192-13	**Date Sampled:**	03/17/17
Matrix:	AQ - Ground Water	**Date Received:**	03/18/17
Method:	SW846-8015C (DAI)	**Percent Solids:**	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109169.D	1	03/27/17	ANJ	n/a	n/a	N:GGH5695
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	72%		56-145%
111-27-3	Hexanol	77%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

SGS Accutest

Report of Analysis

Client Sample ID:	S-37		
Lab Sample ID:	FA42192-14	Date Sampled:	03/17/17
Matrix:	AQ - Ground Water	Date Received:	03/18/17
Method:	SW846-8015C (DAI)	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109170.D	1	03/27/17	ANJ	n/a	n/a	N:GGH5695
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	105%		56-145%
111-27-3	Hexanol	80%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

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ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range
J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

SGS Accutest

Report of Analysis

Client Sample ID:	FB031717		
Lab Sample ID:	FA42192-15	Date Sampled:	03/17/17
Matrix:	AQ - Field Blank Water	Date Received:	03/18/17
Method:	SW846-8015C (DAI)	Percent Solids:	n/a
Project:	BMSMC, Humacao, PR		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1 [a]	GH109171.D	1	03/27/17	ANJ	n/a	n/a	N:GGH5695
Run #2							

Low Molecular Alcohol List

CAS No.	Compound	Result	RL	MDL	Units	Q
64-17-5	Ethanol	ND	100	55	ug/l	
78-83-1	Isobutyl Alcohol	ND	100	36	ug/l	
67-63-0	Isopropyl Alcohol	ND	100	68	ug/l	
71-23-8	n-Propyl Alcohol	ND	100	43	ug/l	
71-36-3	n-Butyl Alcohol	ND	100	87	ug/l	
78-92-2	sec-Butyl Alcohol	ND	100	66	ug/l	
67-56-1	Methanol	ND	200	71	ug/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
111-27-3	Hexanol	75%		56-145%
111-27-3	Hexanol	72%		56-145%

(a) Analysis performed at SGS Accutest, Dayton, NJ.

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

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

FA42192

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PAGE 1 OF 2

Company Name: Anderson Mulholland & Associates
Street Address: 8765 Westchester Avenue, Suite 417
City: Penfield State: NY Zip: 14577
Project Manager: Terry Taylor
Phone #: 914-801-6400
Project Name: 1Qtr Q 2017 Groundwater Sampling - Onsite Wells
Project #:
Client Purchase Order #:
Project Manager: Terry Taylor
Bill to Company Name:
Bill to Street Address:
Bill to City: State: Zip:
Attention:

Matrix Codes: GW - Drinking Water; GW - Ground Water; WW - Water; SW - Surface Water; SO - Soil; SL - Sludge; SED - Sediment; OI - Oil; LIQ - Other Liquid; AIR - Air; SOL - Other Solid; WP - Wipe; PQ - Pool Quality; Equipment Blank; Rinse Blank; Trip Blank

Field ID / Point of Collection	Date	Time	Sampled by	Matrix	# of bottles	VOA	BNA	PEST	HERB	Metals	Cyanide	Sulfide	Phenolics	VOC8260TCL 20	VOC+12XTBE, BROCHL, PTOLU, TBA, TBA, TRF	DRY/GLASS	ASBESTOS	BIOSTORM/VIRUS	BOD-NH3-P-TSS/NAP	SVOC8270TCL, Total Phenols, Total Cyanide, BNAPLUM, CHRYS, DBANTH, H22PYR, NAP	VOC8260TCLM	VMS/PVR	VMS/PVR	VMS/PVR	PESTPESTTCL	LAB USE ONLY
1 PB031617	3/16/17	10:03	NR	WPB	18	9			2	4				X	X	X	X	X	X	X	X	X	X	X	X	
2 S-B-34	3/16/17	11:06	NR	GW	18	9			2	4				X	X	X	X	X	X	X	X	X	X	X	X	
3 S-33	3/16/17	11:19	HR	GW	18	9			2	4				X	X	X	X	X	X	X	X	X	X	X	X	
3 S-33 MS	3/16/17	11:44	HR	WMS	18	9			2	4				X	X	X	X	X	X	X	X	X	X	X	X	
3 S-33 MSD	3/16/17	12:11	HR	WDP	18	9			2	4				X	X	X	X	X	X	X	X	X	X	X	X	
4 TB031617NR	3/16/17	NA	NR	WTB	2	2								X	X						X					
5 TB031617HR	3/16/17	NA	HR	WTB	2	2								X	X						X					
6 S-32	3/16/17	12:00	HR	GW	18	9			2	4				X	X	X	X	X	X	X	X	X	X	X	X	
7 S-36	3/16/17	13:00	HR	GW	18	9			2	4				X	X	X	X	X	X	X	X	X	X	X	X	
8 UP-1	3/17/17	10:00	HR	GW	18	9			2	4				X	X	X	X	X	X	X	X	X	X	X	X	
9 TB031717NR	3/17/17	NA	NR	WTB	2	2								X	X						X					
10 TB031717HR	3/17/17	NA	HR	WTB	2	2								X	X						X					

Turnaround Time (Business days)
- [X] Std. 10 Business Days
- [ ] 5 Day RUSH
- [ ] 4 Day RUSH
- [ ] 3 Day RUSH
- [ ] 1 Day RUSH
- [ ] other

Approved By (MM/Annual PQL) / Date

Data Deliverable Information
- [ ] Commercial "A" (Level 1)
- [ ] Commercial "B" (Level 2)
- [X] FULL71 (Level 3+4)
- [ ] NJ Reduced
- [ ] Commercial "C"
- [ ] NJ Data of Known Quality Protocol Reporting
- [ ] NYSDP Category A
- [ ] NYSDP Category B
- [ ] State Forms
- [ ] EDD Format
- [ ] Other

Commercial "A" = Results Only, Commercial "B" = Results + QC Summary
NJ Reduced = Results + QC Summary + Partial Raw data

Comments / Special Instructions

BANTH, BAPYR, BBFLUAN, BKFLUAN, CHRYS, DBANTH, H22PYR, NAP SHOULD ALL BE ANALYZED BY SIM

Sample Inventory is verified upon client in the Laboratory

Sample Custody must be documented below each time samples change possession, including courier delivery.

Relinquished by	Date/Time	Received by	Relinquished by	Date/Time	Received by
1 [signature]	3/17/17 1500	1 FEDEX	2 FX		1 [signature] 03/18/17
2 Relinquished by Sampler	Date/Time	2	3		2 Relinquished by
3 Relinquished by	Date/Time	3	4		3
4		4			4

Custody Seal # intact / not intact; Preserved where applicable; On Ice; Cooler Temp.

4.8 4.9 4.3 4.6 4.4 1.6 4.7 5.8 5.0

5.1 5

FA42192: Chain of Custody
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FA42192

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PAGE 2 OF 2

Company Name: Anderson Mulholland & Associates
Project Name: 1QT Q 2017 Groundwater Sampling - Onville Wells
Project Manager: Terry Taylor

Lab Sample #	Field ID / Point of Collection	MEOH/H2 Vol #	Date	Time	Sampled By	Matrix	# of Bottles	HCl	NaOH	HNO3	H2SO4	VOSTFCL,B	VOSP-SVTMS, BNZCHL, BTCLLS, TEA, TBA, TBF	DBHELMA	AMSTRL	BNZTRBRHCKX	BNHHMAP,BNHAP	BNPLUMA, CNRTS, DBANTH, H2SPYR, NAP	VMSPSBSTAUN	VMAPVR	VMAPVR	PESTPESTTCL	LAB USE ONLY
11	UP-1DUP		3/17/17	11:08	NR	WDP	10	0		3	4	X	X	X	X	X	X	X	X	X	X	X	
12	S-35		3/17/17	11:12	NR	GW	10	0		3	4	X	X	X	X	X	X	X	X	X	X	X	
13	UP-2		3/17/17	12:39	NR	GW	10	0		3	4	X	X	X	X	X	X	X	X	X	X	X	
14	S-37		3/17/17	12:42	NR	GW	10	0		3	4	X	X	X	X	X	X	X	X	X	X	X	
15	FB0317117		3/17/17	12:10	NR	WFB	10	0		3	4	X	X	X	X	X	X	X	X	X	X	X	
							0																
							0																
							0																
							0																
							0																
							0																
							0																

Matrix Codes: DW - Drinking Water; GW - Ground Water; WW - Water; SW - Surface Water; SO - Soil; SL - Sludge; SED - Sediment; OI - Oil; LIQ - Other Liquid; AIR - Air; SOL - Other Solid; WP - Wipe; FB - Field Blank; EB - Equipment Blank; RB - Rinse Blank; TB - Trip Blank

Turnaround Time (Business days): Approved by (SGS Accutest PM) / Date
- [x] Std. 10 Business Days
- [ ] 4 Day RUSH
- [ ] 3 Day RUSH
- [ ] 2 Day RUSH
- [ ] 1 Day RUSH
- [ ] other

Data Deliverable Information:
- [ ] Commercial "A" (Level 1)
- [ ] Commercial "B" (Level 2)
- [x] FULL1 (Level 3+4)
- [ ] NJ Reduced
- [ ] Commercial "C"
- [ ] All Data of Known Quality Protocol Reporting
- [ ] NYSASP Category A
- [ ] NYSASP Category B
- [ ] State Forms
- [ ] EDD Format
- [ ] Other

Commercial "A" = Results Only, Commercial "B" = Results + QC Summary, NJ Reduced = Results + QC Summary + Partial Raw data

Comments / Special Instructions: BANTH, BAPYRN, BBFLUAN, BKFLUAN, CHRYS, DBANTH, H2SPYR, NAP SHOULD ALL BE ANALYZED BY 8260

Sample Inventory is verified upon receipt in the Laboratory 10/5

Relinquished by Sampler: 1 [signature] 3/17/17 1500
Received by: 1 FEDEX
Relinquished by: 2 Fx
Received by: 2 [signature] 03/18/17
Relinquished by: 3
Received by: 3
Relinquished by: 4
Received by: 4
Custody Seal # 401-400 [ ] Intact [ ] Not Intact
Preserved where applicable [ ]
On Ice [ ]
Cooler Temp.

5.1 5

FA42192: Chain of Custody
Page 2 of 4

EXECUTIVE NARRATIVE

SDG No:	FA42192	Laboratory:	Accutest, Florida
Analysis:	SW846-8260C	Number of Samples:	15
Location:	BMSMC – Humacao, PR		

SUMMARY: Fifteen (15) samples were analyzed for selected volatile organic compounds (VOA Special List) by method SW846-8260C. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence: USEPA Hazardous Waste Support Section SOP No. HW-33A Revision 0 SOM02.2. Low/Medium Volatile Data Validation. July, 2015. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

Critical issues: **None**
Major: **None**
Minor: **None**

Critical findings: **None**
Major findings: **None**
Minor findings:

1. Initial calibration, initial calibration verification, and continuing calibration verification within the method and validation guidance document required performance criteria except for the cases described in the Data Review Worksheet. Closing calibration check verification included in data package.

Results for Carbon tetrachloride and Tetrachloroethene are qualified as estimated (J or UJ) in samples FA42192-13; FA42192-14; and FA42192-15.

Analytes not meeting the continuing calibration verification performance criteria but within the guidance document performance criteria their results are not qualified in affected samples.

2. No target analytes detected in field/trip blanks analyzed except in the cases described in the Data Review Worksheet. Analytes not detected in sample batch above the reporting limits. Results reported as non-detected (U) at the reporting limit.

3. As per chain-of-custody form sample FA42192-3 was to use as the QC sample (MS/MSD) for this batch. Samples FA42192-2 and FA42237-4 were used instead. No action taken, professional judgment.

No MS/MSD associated for batch VP1859 (FA42192-6 diluted); Blank Spike used to assess accuracy. No action taken.

MS/MSD % recoveries and RPD in sample FA42192-2 within laboratory control limits except for the cases described in the Data Review Worksheet.

Results for 1,3-butadiene qualified as estimated (J or UJ) in sample FA42192-2.

Result for tert-butyl alcohol not qualified in affected sample, professional judgment

MS/MSD % recoveries and RPD in sample FA42237-4 within laboratory control limits except for the cases described in the Data Review Worksheet.

No action taken. MS/MSD sample from another job, MS/MSD actions apply only to the field sample used to prepare the Matrix Spike sample.

COMMENTS: Results are valid and can be used for decision making purposes.

Reviewers Name: Rafael Infante
Chemist License 1888

Signature: Rafael Infante

Date: April 26, 2017

SAMPLE ORGANIC DATA SAMPLE SUMMARY

Sample ID: FA42192-1
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: AQ - Field Blank Water

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	1.0	ug/L	1.0	-	U	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Tert-Butyl-Alcohl	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	15.1	ug/L	1.0	-	-	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-2
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	UJ	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	4.1	ug/L	1.0	-	-	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Tert-Butyl Alcohol	31.6	ug/L	1.0	-	-	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-3
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	1.0	ug/L	1.0	-	U	Yes
Bromochloromethane	2.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	0.45	ug/L	1.0	J	J	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	45.1	ug/L	1.0	-	-	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	4.5	ug/L	1.0	-	-	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	6.8	ug/L	1.0	J	J	Yes
Tert-Butyl Alcohol	33.3	ug/L	1.0	-	-	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-4
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: AQ - Trip Blank Water

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	5.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	2.0	ug/L	1.0	-	U	Yes
Carbon disulfide	5.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	1.0	ug/L	1.0	-	U	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Tert-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.8	ug/L	1.0	-	-	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-5
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: AQ - Trip Blank Water

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	1.0	ug/L	1.0	-	U	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	2.7	ug/L	1.0	-	-	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-6
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	5000	ug/L	200	-	U	Yes
Benzene	200	ug/L	200	-	U	Yes
Benzyl Chloride	400	ug/L	200	-	U	Yes
Bromochloromethane	200	ug/L	200	-	U	Yes
Bromodichloromethane	200	ug/L	200	-	U	Yes
Bromoform	200	ug/L	200	-	U	Yes
1,3-Butadiene	400	ug/L	200	-	U	Yes
2-Butanone (MEK)	1000	ug/L	200	-	U	Yes
Carbon disulfide	400	ug/L	200	-	U	Yes
Carbon tetrachloride	200	ug/L	200	-	U	Yes
Chlorobenzene	200	ug/L	200	-	U	Yes
Chloroethane	400	ug/L	200	-	U	Yes
Chloroform	200	ug/L	200	-	U	Yes
Cyclohexane	200	ug/L	200	-	U	Yes
Dibromochloromethane	200	ug/L	200	-	U	Yes
1,2-Dibromo-3-chloropropane	1000	ug/L	200	-	U	Yes
1,2-Dibromoethane	400	ug/L	200	-	U	Yes
Dichlorodifluoromethane	400	ug/L	200	-	U	Yes
1,2-Dichlorobenzene	200	ug/L	200	-	U	Yes
1,3-Dichlorobenzene	200	ug/L	200	-	U	Yes
1,4-Dichlorobenzene	200	ug/L	200	-	U	Yes
1,1-Dichloroethane	200	ug/L	200	-	U	Yes
1,2-Dichloroethane	200	ug/L	200	-	U	Yes
1,1-Dichloroethene	200	ug/L	200	-	U	Yes
cis-1,2-Dichloroethene	200	ug/L	200	-	U	Yes
trans-1,2-Dichloroethene	200	ug/L	200	-	U	Yes

1,2-Dichloropropane	200	ug/L	200	-	U	Yes
cis-1,3-Dichloropropene	200	ug/L	200	-	U	Yes
trans-1,3-Dichloropropene	200	ug/L	200	-	U	Yes
Ethylbenzene	34200	ug/L	2000	-	-	Yes
Freon 113	200	ug/L	200	-	U	Yes
2-Hexanone	2000	ug/L	200	-	U	Yes
Isopropylbenzene	385	ug/L	200	-	-	Yes
p-Isopropyltoluene	200	ug/L	200	-	U	Yes
Methyl Acetate	4000	ug/L	200	-	U	Yes
Methyl Bromide	400	ug/L	200	-	U	Yes
Methyl Chloride	400	ug/L	200	-	U	Yes
Methylcyclohexane	200	ug/L	200	-	U	Yes
Methylene chloride	976	ug/L	200	J	J	Yes
4-Methyl-2-pentanone(MIBK)	1000	ug/L	200	-	U	Yes
Methyl Tert Butyl Ether	200	ug/L	200	-	U	Yes
Styrene	200	ug/L	200	-	U	Yes
Tert-Amyl Alcohol	4000	ug/L	200	-	U	Yes
Ter-Butyl Alcohol	4000	ug/L	200	-	U	Yes
1,1,2,2-Tetrachloroethane	200	ug/L	200	-	U	Yes
Tetrachloroethene	200	ug/L	200	-	U	Yes
Tetrahydrofuran	1000	ug/L	200	-	U	Yes
Toluene	62.7	ug/L	200	J	J	Yes
1,2,3-Trichlorobenzene	400	ug/L	200	-	U	Yes
1,2,4-Trichlorobenzene	400	ug/L	200	-	U	Yes
1,1,1-Trichloroethane	200	ug/L	200	-	U	Yes
1,1,2-Trichloroethane	200	ug/L	200	-	U	Yes
Trichloroethene	200	ug/L	200	-	U	Yes
Trichlorofluoromethane	400	ug/L	200	-	U	Yes
1,2,4-Trimethylbenzene	200	ug/L	200	-	U	Yes
Vinyl chloride	200	ug/L	200	-	U	Yes
m,p-Xylene	56300	ug/L	2000	-	-	Yes
o-Xylene	3650	ug/L	2000	-	-	Yes

Sample ID: FA42192-7
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	8.0	ug/L	1.0	-	-	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	0.61	ug/L	1.0	J	J	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	6.2	ug/L	1.0	J	J	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-8
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	36.9	ug/L	1.0	-	-	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	1.0	ug/L	1.0	-	U	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-9
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: AQ - Trip Blank Water

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	1.0	ug/L	1.0	-	U	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	2.9	ug/L	1.0	-	-	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-10
Sample location: BMSMC, Humacao, PR
Sampling date: 9/23/2016
Matrix: AQ - Trip Blank

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	1.0	ug/L	1.0	-	U	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	3.1	ug/L	1.0	-	-	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-11
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	38.4	ug/L	1.0	-	-	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	1.0	ug/L	1.0	-	U	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-12
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	U	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	1.4	ug/L	1.0	-	-	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	U	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-13
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	UJ	Yes
Chlorobenzene	0.46	ug/L	1.0	J	J	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	2.9	ug/L	1.0	-	-	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	UJ	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-14
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	UJ	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	0.74	ug/L	1.0	J	J	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	UJ	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Sample ID: FA42192-15
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: AQ - Field Blank Water

METHOD: 8260C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Acetone	25	ug/L	1.0	-	U	Yes
Benzene	1.0	ug/L	1.0	-	U	Yes
Benzyl Chloride	2.0	ug/L	1.0	-	U	Yes
Bromochloromethane	1.0	ug/L	1.0	-	U	Yes
Bromodichloromethane	1.0	ug/L	1.0	-	U	Yes
Bromoform	1.0	ug/L	1.0	-	U	Yes
1,3-Butadiene	2.0	ug/L	1.0	-	U	Yes
2-Butanone (MEK)	5.0	ug/L	1.0	-	U	Yes
Carbon disulfide	2.0	ug/L	1.0	-	U	Yes
Carbon tetrachloride	1.0	ug/L	1.0	-	UJ	Yes
Chlorobenzene	1.0	ug/L	1.0	-	U	Yes
Chloroethane	2.0	ug/L	1.0	-	U	Yes
Chloroform	1.0	ug/L	1.0	-	U	Yes
Cyclohexane	1.0	ug/L	1.0	-	U	Yes
Dibromochloromethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dibromo-3-chloropropane	5.0	ug/L	1.0	-	U	Yes
1,2-Dibromoethane	2.0	ug/L	1.0	-	U	Yes
Dichlorodifluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,3-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,4-Dichlorobenzene	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,2-Dichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
cis-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes
trans-1,2-Dichloroethene	1.0	ug/L	1.0	-	U	Yes

1,2-Dichloropropane	1.0	ug/L	1.0	-	U	Yes
cis-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
trans-1,3-Dichloropropene	1.0	ug/L	1.0	-	U	Yes
Ethylbenzene	1.0	ug/L	1.0	-	U	Yes
Freon 113	1.0	ug/L	1.0	-	U	Yes
2-Hexanone	10	ug/L	1.0	-	U	Yes
Isopropylbenzene	1.0	ug/L	1.0	-	U	Yes
p-Isopropyltoluene	1.0	ug/L	1.0	-	U	Yes
Methyl Acetate	20	ug/L	1.0	-	U	Yes
Methyl Bromide	2.0	ug/L	1.0	-	U	Yes
Methyl Chloride	2.0	ug/L	1.0	-	U	Yes
Methylcyclohexane	1.0	ug/L	1.0	-	U	Yes
Methylene chloride	5.0	ug/L	1.0	-	U	Yes
4-Methyl-2-pentanone(MIBK)	5.0	ug/L	1.0	-	U	Yes
Methyl Tert Butyl Ether	1.0	ug/L	1.0	-	U	Yes
Styrene	1.0	ug/L	1.0	-	U	Yes
Tert-Amyl Alcohol	20	ug/L	1.0	-	U	Yes
Ter-Butyl Alcohol	20	ug/L	1.0	-	U	Yes
1,1,2,2-Tetrachloroethane	1.0	ug/L	1.0	-	U	Yes
Tetrachloroethene	1.0	ug/L	1.0	-	UJ	Yes
Tetrahydrofuran	5.0	ug/L	1.0	-	U	Yes
Toluene	1.0	ug/L	1.0	-	U	Yes
1,2,3-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trichlorobenzene	2.0	ug/L	1.0	-	U	Yes
1,1,1-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
1,1,2-Trichloroethane	1.0	ug/L	1.0	-	U	Yes
Trichloroethene	1.0	ug/L	1.0	-	U	Yes
Trichlorofluoromethane	2.0	ug/L	1.0	-	U	Yes
1,2,4-Trimethylbenzene	1.0	ug/L	1.0	-	U	Yes
Vinyl chloride	1.0	ug/L	1.0	-	U	Yes
m,p-Xylene	2.0	ug/L	1.0	-	U	Yes
o-Xylene	1.0	ug/L	1.0	-	U	Yes

Project Number:_FA42192_____
Date:_____March_16-17,_2017_____
Shipping date:__March_17,_2017_____
EPA Region:_____2_____

REVIEW OF VOLATILE ORGANIC PACKAGE
Low/Medium Volatile Data Validation

The following guidelines for evaluating volatile organics were created to delineate required validation actions. This document will assist the reviewer in using professional judgment to make more informed decision and in better serving the needs of the data users. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence: USEPA **Hazardous Waste Support Section SOP No. HW-33A Revision 0 SOM02.2. Low/Medium Volatile Data Validation**. July, 2015. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

The hardcopied (laboratory name) __Accutest_-_Orlando_____ data package received has been reviewed and the quality control and performance data summarized. The data review for VOCs included:

Lab. Project/SDG No.: ____FA42192_____ Sample matrix: __Groundwater___
No. of Samples: _____15_____
Trip blank No.: _____FA42192-4;_FA42192-5;_FA42192-9;_FA42192-10_____
Field blank No.: _____FA42192-1;_ FA42192-15_____
Equipment blank No.:_____-_____
Field duplicate No.:_____-_____

___X___ Data Completeness
___X___ Holding Times
___X___ GC/MS Tuning
___X___ Internal Standard Performance
___X___ Blanks
___X___ Surrogate Recoveries
___X___ Matrix Spike/Matrix Spike Duplicate
___X___ Laboratory Control Spikes
___X___ Field Duplicates
___X___ Calibrations
___X___ Compound Identifications
___X___ Compound Quantitation
___X___ Quantitation Limits

_OverallComments:___Selected_VOA_from_the_special_list_(SW846_8260C)._____
_____

Definition of Qualifiers:

J- Estimated results
U- Compound not detected
R- Rejected data
UJ- Estimated nondetect

Reviewer:_Rafael Infante_____
Date:___April_24,_2017_____

DATA REVIEW WORKSHEETS

DATA COMPLETENESS

MISSING INFORMATION	DATE LAB. CONTACTED	DATE RECEIVED

All criteria were met __X__
Criteria were not met
and/or see below _____

HOLDING TIMES

The objective of this parameter is to ascertain the validity of the results based on the holding time of the sample from time of collection to the time of analysis.

Complete table for all samples and note the analysis and/or preservation not within criteria

SAMPLE ID	DATE SAMPLED	DATE ANALYZED	pH	ACTION
All samples analyzed within method recommended holding. Samples properly preserved.				

Criteria

Aqueous samples – 14 days from sample collection for preserved samples (pH $\leq$ 2, 4$\pm$ 2°C), no air bubbles.
Aqueous samples – 7 days from sample collection for unpreserved samples, 4°C, no air bubbles.
Soil samples- 14 days from sample collection.
Cooler temperature (Criteria: 4 $\pm$ 2 °C): 1.6 – 5.8 °C - OK

Actions

Aqueous samples

a. If there is no evidence that the samples were properly preserved (pH < 2, T = 4°C ± 2°C), but the samples were analyzed within the technical holding time [7 days from sample collection], no qualification of the data is necessary.
b. If there is no evidence that the samples were properly preserved, and the samples were analyzed outside of the technical holding time [7 days from sample collection], qualify detects for all volatile compounds as estimated (J) and non-detects as unusable (R).
c. If the samples were properly preserved, and the samples were analyzed within the technical holding time [14 days from sample collection], no qualification of the data is necessary.
d. If the samples were properly preserved, but were analyzed outside of the technical holding time [14 days from sample collection], qualify detects as estimated (J) and non-detects as unusable (R).
e. If air bubbles were present in the sample vial used for analysis, qualify detected compounds as estimated (J-) and non-detected compounds as estimated (UJ).

Non-aqueous samples

a. If there is no evidence that the samples were properly preserved (T < -7°C or T = 4°C ± 2°C and preserved with $NaHSO_4$), but the samples were analyzed within the technical holding time [14 days

from sample collection], qualify detects for all volatile compounds as estimated (J) and non-detects as (UJ) or unusable (R) using professional judgment.

b. If the samples were properly preserved, and the samples were analyzed within the technical holding time [14 days from sample collection], no qualification of the data is necessary.

c. If there is no evidence that the samples were properly preserved, and the samples were analyzed outside of the technical holding time [14 days from sample collection], qualify detects for all volatile compounds as estimated (J) and non-detects as unusable (R).

d. If the samples were properly preserved, but were analyzed outside of the technical holding time [14 days from sample collection], qualify detects as estimated (J) and non-detects as unusable (R).

Qualify TCLP/SPLP samples

a. If the TCLP/SPLP ZHE procedure is performed within the extraction technical holding time of 14 days, detects and non-detects should not be qualified.

b. If the TCLP/SPLP ZHE procedure is performed outside the extraction technical holding time of 14 days, qualify detects as estimated (J) and non-detects as unusable (R).

c. If TCLP/SPLP aqueous samples and TCLP/SPLP leachate samples are analyzed within the technical holding time of 7 days, detects and non-detects should not be qualified.

d. If TCLP/SPLP aqueous samples and TCLP/SPLP leachate samples are analyzed outside of the technical holding time of 7 days, qualify detects as estimated (J) and non-detects as unusable (R).

Table 1. Holding Time Actions for Low/Medium Volatile Analyses - Summary

Matrix	Preserved	Criteria	Action	
			Detected Associated Compounds	Non-Detected Associated Compounds
Aqueous	No	≤ 7 days	No qualification	
	No	> 7 days	J	R
	Yes	≤ 14 days	No qualification	
	Yes	> 14 days	J	R
Non-Aqueous	No	≤ 14 days	J	Professional judgment. UJ or R
	Yes	≤ 14 days	No qualification	
	Yes No	> 14 days	J	R
TCLP SPLP	Yes	≤ 14 days	No qualification	
TCLP SPLP	No	> 14 days	J	R

TCLP SPLP	ZHE performed within the 14-day technical holding time	No qualification	
TCLP SPLP	ZHE performed outside the 14-day technical holding time	J	R
TCLP SPLP aqueous & TCLP SPLP leachate	Analyzed within 7 days	No qualification	
TCLP SPLP aqueous & TCLP SPLP leachate	Analyzed outside 7 days	J	R
Sample temperature outside 4°C ± 2°C upon receipt at the laboratory		Use professional judgment	
Holding times grossly exceeded		J	R

All criteria were met __X__
Criteria were not met see below _____

GC/MS TUNING

The assessment of the tuning results is to determine if the sample instrumentation is within the standard tuning QC limits

__X___ The BFB performance results were reviewed and found to be within the specified criteria.

__X___ BFB tuning was performed for every 12 hours of sample analysis.

NOTES: All mass spectrometer instrument conditions must be identical to those used during the sample analysis. Background subtraction actions resulting in spectral distortions for the sole purpose of meeting the method specifications are contrary to the Quality Assurance (QA) objectives, and are therefore unacceptable.

NOTES: No data should be qualified based on BFB failure. Instances of this should be noted in the narrative.

All ion abundance ratios must be normalized to m/z 95, the nominal base peak, even though the ion abundance of m/z 174 may be up to 120% that of m/z 95.

Actions:

If samples are analyzed without a preceding valid instrument performance check, qualify all data in those samples as unusable (R).

If ion abundance criteria are not met, professional judgment may be applied to determine to what extent the data may be utilized. When applying professional judgment to this topic, the most important factors to consider are the empirical results that are relatively insensitive to location on the chromatographic profile and the type of instrumentation. Therefore, the critical ion abundance criteria for BFB are the m/z 95/96, 174/175, 174/176, and 176/177 ratios. The relative abundances of m/z 50 and 75 are of lower importance. This issue is more critical for Tentatively Identified Compounds (TICs) than for target analytes.

Note: State in the Data Review Narrative, decisions to use analytical data associated with BFB instrument performance checks not meeting contract requirements.

Note: Verify that that instrument instrument performance check criteria were achieved using techniques described in Low/Medium Volatiles Organic Analysis, Section II.D.5 of the SOM02.2 NFG, obtain additional information on the instrument performance checks. Make sure that background subtraction was performed from the BFB peak and not from background subtracting from the solvent front or from another region of the chromatogram.

Use professional judgment to determine whether associated data should be qualified based on the spectrum of the mass calibration compound.

List the samples affected:

_____
_____
_____
_____
_____
_____
_____

If mass calibration is in error, all associated data are rejected.

All criteria were met _____
Criteria were not met
and/or see below ___X___

CALIBRATION VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:_____03/28/17_____	_____03/28/17_____
Dates of continuing (initial) calibration:__03/28/17_____	_____03/28/17_____
Dates of continuing calibration:_____03/29/17_____	_____03/29/17;_03/30/17_____
Dates of ending calibration:___03/28/17;_03/29/17_____	__03/28/17;_03/29/17;_03/30/17_
Instrument ID numbers:_____GCMSM_____	_____GCMSP_____
Matrix/Level:_____Aqueous/low_____	_____Aqueous/low_____

DATE	LAB FILE ID#	CRITERIA OUT RFs, %RSD, **%D**, r	COMPOUND	SAMPLES AFFECTED
GCMSM				
03/28/17	icv4065-5	20.8	Bromomethane^	FA42192-13 to 15
		25.5	Acetone^	
03/29/17	CC4065-5	- 29.5	Carbon tetrachloride ✓	FA42192-13 to -15
		- 25.5	Tetrachloroethene ✓	
		20.4	1,2-dibromo-3-chloropropane^	
		-21.4	1,2,3-trichlorobenzene^	
GCMSP				
03/28/17	ICV1854-5	21.4	Chloroethane^	FA42192-1 to -12

Note: Initial calibration, initial calibration verification, and continuing calibration verification within the method and validation guidance document required performance criteria except for the cases described in this document. Closing calibration check verification included in data package.

Results for Carbon tetrachloride and Tetrachloroethene are qualified as estimated (J or UJ) in samples FA42192-13; FA42192-14; and FA42192-15.

^Analyte not meeting the continuing calibration verification performance criteria but within the guidance document performance criteria. Results are not qualified in affected samples.

Criteria

The analyte calibration criteria in the following Table must be obtained. Analytes not meeting the criteria are qualified.

A separate worksheet should be filled for each initial curve.

Initial Calibration - Table 2. RRF, %RSD, and %D Acceptance Criteria for Initial Calibration and CCV for Low/Medium Volatile Analysis

Analyte	Minimum RRF	Maximum %RSD	Opening Maximum %D[1]	Closing Maximum %D
Dichlorodifluoromethane	0.010	25.0	±40.0	±50.0
Chloromethane	0.010	20.0	±30.0	±50.0
Vinyl chloride	0.010	20.0	±25.0	±50.0
Bromomethane	0.010	40.0	±30.0	±50.0
Chloroethane	0.010	40.0	±25.0	±50.0
Trichlorofluoromethane	0.010	40.0	±30.0	±50.0
1.1-Dichloroethene	0.060	20.0	±20.0	±25.0
1.1.2-Trichloro-1.2.2-trifluoroethane	0.050	25.0	±25.0	±50.0
Acetone	0.010	40.0	±40.0	±50.0
Carbon disulfide	0.100	20.0	±25.0	±25.0
Methyl acetate	0.010	40.0	±40.0	±50.0
Methylene chloride	0.010	40.0	±30.0	±50.0
trans-1.2-Dichloroethene	0.100	20.0	±20.0	±25.0
Methyl tert-butyl ether	0.100	40.0	±25.0	±50.0
1.1-Dichloroethane	0.300	20.0	±20.0	±25.0
cis-1.2-Dichloroethene	0.200	20.0	±20.0	±25.0
2-Butanone	0.010	40.0	±40.0	±50.0
Bromochloromethane	0.100	20.0	±20.0	±25.0
Chloroform	0.300	20.0	±20.0	±25.0
1.1.1-Trichloroethane	0.050	20.0	±25.0	±25.0
Cyclohexane	0.010	40.0	±25.0	±50.0
Carbon tetrachloride	0.100	20.0	±25.0	±25.0
Benzene	0.200	20.0	±20.0	±25.0
1.2-Dichloroethane	0.070	20.0	±20.0	±25.0
Trichloroethene	0.200	20.0	±20.0	±25.0
Methylcyclohexane	0.050	40.0	±25.0	±50.0
1.2-Dichloropropane	0.200	20.0	±20.0	±25.0
Bromodichloromethane	0.300	20.0	±20.0	±25.0
cis-1.3-Dichloropropene	0.300	20.0	±20.0	±25.0
4-Methyl-2-pentanone	0.030	25.0	±30.0	±50.0
Toluene	0.300	20.0	±20.0	±25.0
trans-1.3-Dichloropropene	0.200	20.0	±20.0	±25.0
1.1.2-Trichloroethane	0.200	20.0	±20.0	±25.0
Tetrachloroethene	0.100	20.0	±20.0	±25.0
2-Hexanone	0.010	40.0	±40.0	±50.0
Dibromochloromethane	0.200	20.0	±20.0	±25.0
1.2-Dibromoethane	0.200	20.0	±20.0	±25.0
Chlorobenzene	0.400	20.0	±20.0	±25.0
Ethylbenzene	0.400	20.0	±20.0	±25.0

Analyte	Minimum RRF	Maximum %RSD	Opening Maximum %D[1]	Closing Maximum
m,p-Xylene	0.200	20.0	±20.0	±25.0
o-Xylene	0.200	20.0	±20.0	±25.0
Styrene	0.200	20.0	±20.0	±25.0
Bromoform	0.100	20.0	±25.0	±50.0
Isopropylbenzene	0.400	20.0	±25.0	±25.0
1,1,2,2-Tetrachloroethane	0.200	20.0	±25.0	±25.0
1,3-Dichlorobenzene	0.500	20.0	±20.0	±25.0
1,4-Dichlorobenzene	0.600	20.0	±20.0	±25.0
1,2-Dichlorobenzene	0.600	20.0	±20.0	±25.0
1,2-Dibromo-3-chloropropane	0.010	25.0	±30.0	±50.0
1,2,4-Trichlorobenzene	0.400	20.0	±30.0	±50.0
1,2,3-Trichlorobenzene	0.400	25.0	±30.0	±50.0
Deuterated Monitoring Compound				
Vinyl chloride-d_3	0.010	20.0	±30.0	±50.0
Chloroethane-d_5	0.010	40.0	±30.0	±50.0
1,1-Dichloroethene-d_2	0.050	20.0	±25.0	±25.0
2-Butanone-d_5	0.010	40.0	±40.0	±50.0
Chloroform-d	0.300	20.0	±20.0	±25.0
1,2-Dichloroethane-d_4	0.060	20.0	±25.0	±25.0
Benzene-d_6	0.300	20.0	±20.0	±25.0
1,2-Dichloropropane-d_6	0.200	20.0	±20.0	±25.0
Toluene-d_8	0.300	20.0	±20.0	±25.0
trans-1,3-Dichloropropene-d_4	0.200	20.0	±20.0	±25.0
2-Hexanone-d_5	0.010	40.0	±40.0	±50.0
1,1,2,2-Tetrachloroethane-d_2	0.200	20.0	±25.0	±25.0
1,2-Dichlorobenzene-d_4	0.400	20.0	±20.0	±25.0

[1] If a closing CCV is acting as an opening CCV, all target analytes and DMCs must meet the requirements for an opening CCV.

Actions:

1. If any volatile target compound has an RRF value less than the minimum in the table, use professional judgment for detects, based on mass spectral identification, to qualify the data as estimated (J+ or R).
   a. If any volatile target compound has an RRF value less than the minimum criterion, qualify non-detected compounds as unusable (R).
   b. If any of the volatile target compounds listed in the Table has %RSD greater than the criteria, qualify detects as estimated (J), and non-detected compounds using professional judgment.
   c. If the volatile target compounds meet the acceptance criteria for RRF and the %RSD, no qualification of the data is necessary.

d. No qualification of the data is necessary on the DMC RRF and %RSD data alone. Use professional judgment and follow the guidelines in Action 2 to evaluate the DMC RRF and %RSD data in conjunction with the DMC recoveries to determine the need for qualification of data.

2. At the reviewer's discretion, and based on the project-specific Data Quality Objectives (DQOs), a more in-depth review may be considered using the following guidelines:
   a. If any volatile target compound has a %RSD greater than the maximum criterion in the Table, and if eliminating either the high or the low-point of the curve does not restore the %RSD to less than or equal to the required maximum:
      i. Qualify detects for that compound(s) as estimated (J).
      ii. Qualify non-detected volatile target compounds using professional judgment.
   b. If the high-point of the curve is outside of the linearity criteria (e.g., due to saturation):
      i. Qualify detects outside of the linear portion of the curve as estimated (J).
      ii. No qualifiers are required for detects in the linear portion of the curve.
      iii. No qualifiers are required for volatile target compounds that were not detected.
   c. If the low-point of the curve is outside of the linearity criteria:
      i. Qualify low-level detects in the area of non-linearity as estimated (J).
      ii. No qualifiers are required for detects in the linear portion of the curve.
      iii. For non-detected volatile compounds, use the lowest point of the linear portion of the curve to determine the new quantitation limit.

Note: If the laboratory has failed to provide adequate calibration information, inform the Region's designated representative to contact the laboratory and request the necessary information. If the information is not available, the reviewer must use professional judgment to assess the data.

State in the Data Review Narrative, if possible, the potential effects on the data due to calibration criteria exceedance.

Note, for the Laboratory COR action, if calibration criteria are grossly exceeded.

Table. Initial Calibration Actions for Low/Medium Volatile Analysis – Summary

Criteria	Action	
	Detect	Non-detect
Initial Calibration not performed at specified frequency and sequence	Use professional judgment R	Use professional judgment R
Initial Calibration not performed at the specified concentrations	J	UJ
RRF < Minimum RRF in Table for target analyte	Use professional judgment J+ or R	R
RRF > Minimum RRF in Table for target analyte	No qualification	No qualification
%RSD > Maximum %RSD in Table for target analyte	J	Use professional judgment
%RSD ≤ Maximum %RSD in Table for target analyte	No qualification	No qualification

All criteria were met __X___
Criteria were not met
and/or see below _____

Continuing Calibration Verification (CCV)

NOTE: Verify that the CCV was run at the required frequency (an opening and closing CCV must be run within 12-hour period) and the CCV was compared to the correct initial calibration. If the mid-point standard from the initial calibration is used as an opening CCV, verify that the result (RRF) of the mid-point standard was compared to the average RRF from the correct initial calibration.

The closing CCV used to bracket the end of a 12-hour analytical sequence may be used as the opening CCV for the new 12-hour analytical sequence, provided that all the technical acceptance criteria are met for an opening CCV (see criteria show before in the Table) . If the closing CCV does not meet the technical acceptance criteria for an opening CCV, then a BFB tune followed by an opening CCV is required and the next 12-hour time period begins with the BFB tune.

All DMCs must meet RRF criteria. No qualification of the data is necessary on the DMCs RRF and %RSD/%D data alone. However, use professional judgment to evaluate the DMC and %RSD/%D data in conjunction with the DMC recoveries to determine the need of qualification the data.

Action:

1. If a CCV (opening and closing) was not run at the appropriate frequency, qualify data using professional judgment.

2. Qualify all volatile target compounds in Table shown before using the following criteria:

   a. For an opening CCV, if any volatile target compound has an RRF value less than the minimum criterion, use professional judgment for detects, based on mass spectral identification, to qualify the data as estimated (J) and qualify non-detected compounds as unusable (R).
   b. For a closing CCV, if any volatile target compound has an RRF value less than the criteria, use professional judgment for detects based on mass spectral identification to qualify the data as estimated (J), and qualify non-detected compounds as unusable (R).
   c. For an opening CCV, if the Percent Difference value for any of the volatile target compounds is outside the limits in calibration criteria Table shown before, qualify detects as estimated (J) and non-detected compounds as estimated (UJ).
   d. For a closing CCV, if the Percent Difference value for any volatile target compound is outside the limits in calibration criteria table, qualify detects as estimated (J) and non-detected compounds as estimated (UJ).
   e. If the volatile target compounds meet the acceptable criteria for RRF and the Percent Difference, no qualification of the data is necessary.

f. No qualification of the data is necessary on the DMC RRF and the Percent Difference data alone. Use professional judgment to evaluate the DMC RRF and Percent Difference data in conjunction with the DMC recoveries to determine the need for qualification of data.

Notes: If the laboratory has failed to provide adequate calibration information, inform the Region's designated representative to contact the laboratory and request the necessary information. If the information is not available, the reviewer must use professional judgment to assess the data.

State in the Data Review Narrative, if possible, the potential effects on the data due to calibration criteria exceedance.

Note, for Contract Laboratory COR action, if calibration criteria are grossly exceeded.

Table. Continuing Calibration Actions for Low/Medium Volatile Analysis – Summary

Criteria for Opening CCV	Criteria for Closing CCV	Action	
		Detect	Non-detect
CCV not performed at required frequency	CCV not performed at required frequency	Use professional judgment R	Use professional judgment R
CCV not performed at specified concentration	CCV not performed at specified concentration	Use professional judgment	Use professional judgment
RRF < Minimum RRF in Table 2 for target analyte	RRF < Minimum RRF in Table for target analyte	Use professional judgment J or R	R
RRF > Minimum RRF in Table 2 for target analyte	RRF > Minimum RRF in Table for target analyte	No qualification	No qualification
%D outside the Opening Maximum %D limits in Table 2 for target analyte	%D outside the Closing Maximum %D limits in Table for target analyte	J	UJ
%D within the inclusive Opening Maximum %D limits in Table 2 for target analyte	%D within the inclusive Closing Maximum %D limits in Table for target analyte	No qualification	No qualification

All criteria were met ___X___
Criteria were not met
and/or see below _____

BLANK ANALYSIS RESULTS (Sections 1 & 2)

The assessment of the blank analysis results is to determine the existence and magnitude of contamination problems. The criteria for evaluation of blanks apply only to blanks associated with the samples, including trip, equipment, and laboratory blanks. If problems with any blanks exist, all data associated with the case must be carefully evaluated to determine whether or not there is an inherent variability in the data for the case, or if the problem is an isolated occurrence not affecting other data.

List the contamination in the blanks below. High and low levels blanks must be treated separately.

The concentration of a target analyte in any blank must not exceed its Contract Required Quantitation Limit (CRQL) (2x CRQLs for Methylene chloride, Acetone, and 2-Butanone). TIC concentration in any blanks must be ≤ 5.0 µg/L for water (0.0050 mg/L for TCLP leachate) and ≤ 5.0 µg/kg for soil matrices.

Laboratory blanks

The method blank, like any other sample in the SDG, must meet the technical acceptance criteria for sample analysis.

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
_No_target_analyte_detected_in_method_blanks._				

Field/Equipment/**Trip blank**

If field or trip blanks are present, the data reviewer should evaluate this data in a similar fashion as the method blanks.

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
_No_target_analytes_detected_in_field/trip_blanks_analyzed_except_in_the_cases_described_in__ _this_document._No_equipmen_blank_analyzed_in_this_sample_batch._____				
_03/29/17_____	FA42192-1___	Aqueous/low_____	Trichloroethylene_____	15.1_ug/l_____
_03/29/17_____	FA42192-4___	Aqueous/low_____	Toluene_____	1.8__ug/l_____
_03/29/17_____	FA42192-5___	Aqueous/low_____	Toluene_____	2.7__ug/l_____
_03/29/17_____	FA42192-9___	Aqueous/low_____	Toluene_____	2.9__ug/l_____
_03/29/17_____	FA42192-10__	Aqueous/low_____	Toluene_____	3.1__ug/l_____

Note: Analytes not detected in sample batch above the reporting limits. Results reported as non-detected (U) at the reporting limit.

All criteria were met __X___
Criteria were not met
and/or see below _____

BLANK ANALYSIS RESULTS (Section 3)

Blank Actions

Note: All fields blank results associated with a particular group of samples (may exceed one per case) must be used to qualify data. Trip blanks are used to qualify only those samples with which they were shipped. Blanks may not be qualified because of contamination in another blank. Field blanks and trip blanks must be qualified for system monitoring compounds, instrument performance criteria, and spectral or calibration QC problems.

Samples taken from a drinking water tap do not have associated field blanks.

When applied as described in the Table below, the contaminant concentration in the blank is multiplied by the sample dilution factor.

Table. Blank and TCLP/SPLP LEB Actions for Low/Medium Volatile Analysis

Blank Type	Blank Result	Sample Result	Action for Samples
Method. Storage. Field. Trip. TCLP SPLP LEB. Instrument**	Detects	Not detected	No qualification required
	< CRQL *	< CRQL*	Report CRQL value with a U
		≥ CRQL*	No qualification required
	> CRQL *	< CRQL*	Report CRQL value with a U
		≥ CRQL* and ≤ blank concentration	Report blank value for sample concentration with a U
		≥ CRQL* and > blank concentration	No qualification required
	= CRQL*	≤ CRQL*	Report CRQL value with a U
		> CRQL*	No qualification required
	Gross contamination	Detects	Report blank value for sample concentration with a U

* 2x the CRQL for methylene chloride, 2-butanone and acetone.
** Qualifications based on instrument blank results affect only the sample analyzed immediately after the sample that has target compounds that exceed the calibration range or non-target compounds that exceed 100 µg/L.

Action Levels (ALs) should be based upon the highest concentration of contaminant determined in any blank. Do not qualify any blank with another blank. The ALs for samples which have been diluted should be corrected for the sample dilution factor and/or % moisture, where applicable. No positive sample results should be reported unless the concentration of the compound in the samples exceeds the ALs:

Notes:

High and low level blanks must be treated separately
Compounds qualified "U" for blank contamination are still considered "hits" when qualifying for calibration criteria.

CONTAMINATION SOURCE/LEVEL	COMPOUND	CONC/UNITS	AL/UNITS	SQL	AFFECTED SAMPLES

All criteria were met __X__
Criteria were not met
and/or see below _____

DEUTERATED MONITORING COMPOUNDS (DMCs)

Laboratory performance of individual samples is established by evaluation of surrogate spike (DMCs) recoveries. All samples are spiked with surrogate compounds prior to sample analysis. The accuracy of the analysis is measured by the surrogate percent recovery. Since the effects of the sample matrix are frequently outside the control of the laboratory and may present relatively unique problems, the validation of data is frequently subjective and demands analytical experience and professional judgment.

Table. Volatile Deuterated Monitoring Compounds (DMCs) and Recovery Limits

DMC	%R for Water Sample	%R for Soil Sample
Vinyl chloride-d3	60-135	30-150
Chloroethane-d5	70-130	30-150
1.1-Dichloroethene-d2	60-125	45-110
2-Butanone-d5	40-130	20-135
Chloroform-d	70-125	40-150
1.2-Dichloroethane-d4	70-125	70-130
Benzene-d6	70-125	20-135
1.2-Dichloropropane-d6	70-120	70-120
Toluene-d8	80-120	30-130
trans-1.3-Dichloropropene-d4	60-125	30-135
2-Hexanone-d5	45-130	20-135
1.1.2.2-Tetrachloroethane-d2	65-120	45-120
1.2-Dichlorobenzene-d4	80-120	75-120

NOTE: The recovery limits for any of the compounds listed in the above Table may be expanded at any time during the period of performance if the United States Environmental Protection Agency (EPA) determines that the limits are too restrictive.

Action:

Are recoveries for DMCs in volatile samples and blanks must be within the limits specified in the Table above. **Yes**? or No?

NOTE: The recovery limits for any of the compounds listed in the Table above may be expanded at any time during the period of performance if USEPA determines that the limits are too restrictive.

List the DMCs that may fail to meet the recovery limits

Sample ID	Date	DMCs	% Recovery	Action

Note: DMCs recoveries within the laboratory required control limits and within the guidance document performance criteria (80 – 120). Other non-deuterated surrogates added to the samples within laboratory control limits.

Note: Any sample which has more than 3 DMCs outside the limits must be reanalyzed.

Action:

1. For any recovery greater than the upper acceptance limit:
   a. Qualify detected associated volatile target compounds as estimated high (J+).
   b. Do not qualify non-detected associated volatile target compounds.
2. For any recovery greater than or equal to 10%, and less than the lower acceptance limit:
   a. Qualify detected associated volatile target compounds as estimated low (J-).
   b. Qualify non-detected associated volatile target compounds as estimated (UJ).
3. For any recovery less than 10%:
   a. Qualify detected associated volatile target compounds as estimated low (J-).
   b. Qualify non-detected associated volatile target compounds as unusable (R).
4. For any recovery within acceptance limits, no qualification of the data is necessary.
5. In the special case of a blank analysis having DMCs out of specification, the reviewer must give special consideration to the validity of associated sample data. The basic concern is whether the blank problems represent an isolated problem with the blank alone, or whether there is a fundamental problem with the analytical process. For example, if one or more samples in the batch show acceptable DMC recoveries, the reviewer may choose to consider the blank problem to be an isolated occurrence. However, even if this judgment allows some use of the affected data, note analytical problems for Contract Laboratory COR action.
6. If more than three DMCs are outside of the recovery limits for Low/Medium volatiles analysis and the sample was not reanalyzed, note under Contract Problems/Non-Compliance.

Table. Deuterated Monitoring Compound (DMC) Recovery Actions for Low/Medium Volatiles Analyses – Summary

Criteria	Action	
	Detect Associated Compounds	Non-detected Associated Compounds
%R < 10%	J-	R
10% ≤ %R < Lower Acceptance Limit	J-	UJ
Lower Acceptance Limit ≤ %R ≤ Upper Acceptance Limit	No qualification	No qualification
%R > Upper Acceptance Limit	J+	No qualification

TABLE. VOLATILE DEUTERATED MONITORING COMPOUNDS (DMCs) AND THE ASSOCIATED TARGET COMPOUNDS

Vinyl chloride-d_3 (DMC-1)	**Chloroethane-d_5 (DMC-2)**	**1,1-Dichloroethene-d_2 (DMC-3)**
Vinyl chloride	Dichlorodifluoromethane Chloromethane Bromomethane Chloroethane Carbon disulfide	trans-1,2-Dichloroethene cis-1,2-Dichloroethene 1,1-Dichloroethene
2-Butanone-d_5 (DMC-4)	**Chloroform-d (DMC-5)**	**1,2-Dichloroethane-d_4 (DMC-6)**
Acetone 2-Butanone	1,1-Dichloroethane Bromochloromethane Chloroform Dibromochloromethane Bromoform	Trichlorofluoromethane 1,1,2-Trichloro-1,2,2-trifluoroethane Methyl acetate Methylene chloride Methyl-tert-butyl ether 1,1,1-Trichloroethane Carbon tetrachloride 1,2-Dibromoethane 1,2-Dichloroethane
Benzene-d_6 (DMC-7)	**1,2-Dichloropropane-d_6 (DMC-8)**	**Toluene-d_8 (DMC-9)**
Benzene	Cyclohexane Methylcyclohexane 1,2-Dichloropropane Bromodichloromethane	Trichloroethene Toluene Tetrachloroethene Ethylbenzene o-Xylene m,p-Xylene Styrene Isopropylbenzene
trans-1,3-Dichloropropene-d_4 (DMC-10)	**2-Hexanone-d_5 (DMC-11)**	**1,1,2,2-Tetrachloroethane-d_2 (DMC-12)**
cis-1,3-Dichloropropene trans-1,3-Dichloropropene 1,1,2-Trichloroethane	4-Methyl-2-pentanone 2-Hexanone	1,1,2,2,-Tetrachloroethane 1,2-Dibromo-3-chloropropane
1,2-Dichlorobenzene-d_4 (DMC-13)		
Chlorobenzene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene 1,2,4-Trichlorobenzene 1,2,3-Trichlorobenzene		

All criteria were met ___X_____
Criteria were not met
and/or see below _____

MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

This data is generated to determine long term precision and accuracy in the analytical method for various matrices. This data alone cannot be used to evaluate the precision and accuracy of individual samples. If any % R in the MS or MSD falls outside the designated range, the reviewer should determine if there are matrix effects, i.e. LCS data are within the QC limits but MS/MSD data are outside QC limit.

NOTES: Data for MS and MSDs will not be present unless requested by the Region. Notify the Contract Laboratory COR if a field or trip blank was used for the MS and MSD.

For a Matrix Spike that does not meet criteria, apply the action to only the field sample used to prepare the Matrix Spike sample. If it is clearly stated in the data validation materials that the samples were taken through incremental sampling or some other method guaranteeing the homogeneity of the sample group, then the entire sample group may be qualified.

1. MS/MSD Recoveries and Precision Criteria

The laboratory should use one MS and a duplicate analysis of an unspiked field sample if target analytes are expected in the sample. If target analytes are not expected, MS/MSD should be analyzed.

List the %Rs, RPD of the compounds which do not meet the criteria.

Sample ID:_ FA42192-2MS/-2MSD____ Matrix/Level:__Groundwater___
Sample ID:_ FA42237-4MS/-4MSD____ Matrix/Level:__Groundwater___

The QC reported here applies to the following samples: Method: **SW846 8260C**
FA42192-1, FA42192-2, FA42192-3; FA42192-4, FA42192-5, FA42192-6; FA42192-7, FA42192-8, FA42192-9; FA42192-10; FA42192-11; FA42192-12

Compound	FA42192-2 ug/l	Q	Spike ug/l	MS ug/	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
1,3-Butadiene ✓	ND		25	9.0	36*	25	9.3	37*	3	65-135/31
Tert-Butyl Alcohol ✓	31.6		250	352	128	250	367	134*	4	63-129/27

* Outside laboratory control limits

Note: As per chain-of-custody form sample FA42192-3 was to use as the QC sample (MS/MSD) for this batch. Samples FA42192-2 and FA42237-4 were used instead. No action taken, professional judgment.

No MS/MSD associated for batch VP1859 (FA42192-6 diluted); Blank Spike used to assess accuracy. No action taken.

MS/MSD % recoveries and RPD within laboratory control limits except for the cases described in this document.

Results for 1,3-butadiene qualified as estimated (J or UJ) in sample FA42192-2.

Result for tert-butyl alcohol not qualified in affected sample, professional judgment

The QC reported here applies to the following samples: Method: **SW846 8260C**
FA42192-13; FA42192-14, FA42192-15

Compound	FA42237-4 ug/l	Q	Spike ug/l	MS ug/	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
1,3-Butadiene ✓	ND		25	1.7	7*	25	1.6	6*	6	65-135/31

Note: MS/MSD % recoveries and RPD within laboratory control limits except for the cases described in this document.

No action taken. MS/MSD sample from another job, MS/MSD actions apply only to the field sample used to prepare the Matrix Spike sample.

Note:

* QC limits are laboratory in-house performance criteria, LL = lower limit, UL = upper limit.
* If QC limits are not available, use limits of 70 – 130 %.

Actions:

1. No qualification of the data is necessary on MS and MSD data alone. However, using professional judgment, the validator may use the MS and MSD results in conjunction with other QC criteria and determine the need for some qualification of the data.

QUALITY	%R < LL	%R > UL
Positive results	J	J
Nondetects results	R	Accept

MS/MSD criteria apply only to the unspiked sample, its dilutions, and the associated MS/MSD samples:

If the % R for the affected compounds were < LL (or 70 %), qualify positive results (J) and nondetects (UJ).
If the % R for the affected compounds were > UL (or 130 %), only qualify positive results (J).
If 25 % or more of all MS/MSD %R were < LL (or 70 %) or if two or more MS/MSD %Rs were < 10%, qualify all positive results (J) and reject nondetects (R).

A separate worksheet should be used for each MS/MSD pair.

All criteria were met __X____
Criteria were not met
and/or see below _____

LABORATORY CONTROL SAMPLE (LCS) ANALYSIS

This data is generated to determine accuracy of the analytical method for various matrices.

1. LCS Recoveries Criteria

Where LCS spiked with the same analyte at the same concentrations as the MS/MSD? **Yes** or No. If no make note in data review memo.
List the %R of compounds which do not meet the criteria

LCS ID COMPOUND % R QC LIMIT

Recoveries(blank_spike)_within_laboratory_control_limits._____

_____

_____

_____

_____

_____

Note:

* QC limits are laboratory in-house performance criteria, LL = lower limit, UL = upper limit.
* If QC limits are not available, use limits of 70 – 130 %.

Actions:

QUALITY	%R < LL	%R > UL
Positive results	J	J
Nondetects results	R	Accept

All analytes in the associated sample results are qualified for the following criteria.

If 25 % of the LCS recoveries were < LL (or 70 %), qualify all positive results (j) and reject nondetects (R).
If two or more LCS were below 10 %, qualify all positive results as (J) and reject nondetects (R).

2. Frequency Criteria:

Where LCS analyzed at the required frequency and for each matrix? **Yes** or No.
If no, the data may be affected. Use professional judgment to determine the severity of the effect and qualify data accordingly. Discuss any actions below and list the samples affected.

All criteria were met ___X___
Criteria were not met
and/or see below _____

IX. FIELD/LABORATORY DUPLICATE PRECISION

Sample IDs: __FA42192-8/-11_____ Matrix:__Groundwater___

Field/laboratory duplicates samples may be taken and analyzed as an indication of overall precision. These analyses measure both field and lab precision; therefore, the results may have more variability than laboratory duplicates which only laboratory performance. It is also expected that soil duplicate results will have a greater variance than water matrices due to difficulties associated with collecting identical field duplicate samples.

The project QAPP should be reviewed for project-specific information.

> **NOTE:** In the absence of QAPP guidance for validating data from field duplicates, the following action will be taken.

Identify which samples within the data package are field duplicates. Estimate the relative percent difference (RPD) between the values for each compound. Use professional judgment to note large RPDs (> 50%) in the narrative.

COMPOUND	SQL	SAMPLE CONC.	DUPLICATE CONC.	RPD	ACTION
Field duplicate analyzed with this data package. RPD within required criteria, ≤ 50 % for target analytes detected at concentration > 5x the SQL.					

Actions:

Qualify as estimated positive results (J) and nondetects (UJ) for the compound that exceeded the above criteria. For organics, only the sample and duplicate will be qualified.

If an RPD cannot be calculated because one or both of the sample results is not detected, the following actions are suggested based on professional judgment:

If one sample result is not detected and the other is greater than 5x the SQL qualify (J/UJ).

If one sample value is not detected and the other is greater than 5x the SQL and the SQLs for the sample and duplicate are significantly different, use professional judgment to determine if qualification is appropriate.

If one sample value is not detected and the other is less than 5x, use professional judgment to determine if qualification is appropriate.

If both sample and duplicate results are not detected, no action is needed.

All criteria were met ___X___
Criteria were not met
and/or see below _____

X. INTERNAL STANDARD PERFORMANCE

The assessment of the internal standard (IS) parameter is used to assist the data reviewer in determining the condition of the analytical instrumentation.

DATE	SAMPLE ID	IS OUT	IS AREA	ACCEPTABLE RANGE	ACTION
Internal standard area within laboratory control limits.					

Action:

1. If an internal standard area count for a sample or blank is greater than 200.0% of the area for the associated standard (opening CCV or mid-point standard from initial calibration) (see Table below):
   a. Qualify detects for compounds quantitated using that internal standard as estimated low (J-).
   b. Do not qualify non-detected associated compounds.
2. If an internal standard area count for a sample or blank is less than 20.0% of the area for the associated standard (opening CCV or mid-point standard from initial calibration):
   a. Qualify detects for compounds quantitated using that internal standard as estimated high (J+).
   b. Qualify non-detected associated compounds as unusable (R).
3. If an internal standard area count for a sample or blank is greater than or equal to 20.0%, and less than or equal to 200% of the area for the associated standard opening CCV or mid-point standard from initial calibration, no qualification of the data is necessary.
4. If an internal standard RT varies by more than 30.0 seconds: Examine the chromatographic profile for that sample to determine if any false positives or negatives exist. For shifts of a large magnitude, the reviewer may consider partial or total rejection of the data for that sample fraction. Detects should not need to be qualified as unusable (R) if the mass spectral criteria are met.
5. If an internal standard RT varies by less than or equal to 30.0 seconds, no qualification of the data is necessary.

Note: Inform the Contract Laboratory Program Project Officer (CLP PO) if the internal standard performance criteria are grossly exceeded. Note in the Data Review Narrative potential effects on the data resulting from unacceptable internal standard performance.

6. If required internal standard compounds are not added to a sample or blank, qualify detects and non-detects as unusable (R).
7. If the required internal standard compound is not analyzed at the specified concentration in a sample or blank, use professional judgment to qualify detects and non-detects.

Table. Internal Standard Actions for Low/Medium Volatiles Analyses - Summary

Criteria	Action	
	Detected Associated Compounds*	Non-detected Associated Compounds*
Area counts > 200% of 12-hour standard (opening CCV or mid-point standard from initial calibration)	J-	No qualification
Area counts < 20% of 12-hour standard (opening CCV or mid-point standard from initial calibration)	J+	R
Area counts ≥ 50% but ≤ 200% of 12-hour standard (opening CCV or mid-point standard from initial calibration)	No qualification	
RT difference > 30.0 seconds between samples and 12-hour standard (opening CCV or mid-point standard from initial calibration)	R **	R
RT difference ≤ 30.0 seconds between samples and 12-hour standard (opening CCV or mid-point standard from initial calibration)	No qualification	

* For volatile compounds associated to each internal standard, see TABLE - VOLATILE TARGET ANALYTES, DEUTERATED MONITORING COMPOUNDS WITH ASSOCIATED INTERNAL STANDARDS FOR QUANTITATION in SOM02.2, Exhibit D, available at: http://www.epa.gov/superfund/programs/clp/download/som/som22d.pdf
** Detects should not need to be qualified as unusable (R) if the mass spectral criteria are met.

All criteria were met _X__
Criteria were not met
and/or see below _____

TARGET COMPOUND IDENTIFICATION

Criteria:

Is the Relative Retention Times (RRTs) of reported compounds within ±0.06 RRT units of the standard RRT [opening Continuing Calibration Verification (CCV) or mid-point standard from the initial calibration]. **Yes**? or No?

List compounds not meeting the criteria described above:

Sample ID	Compounds	Actions

Mass spectra of the sample compound and a current laboratory-generated standard [i.e., the mass spectrum from the associated calibration standard (opening CCV or mid-point standard from initial calibration)] must match according to the following criteria:

a. All ions present in the standard mass spectrum at a relative intensity greater than 10% must be present in the sample spectrum.
b. The relative intensities of these ions must agree within ±20% between the standard and sample spectra (e.g., for an ion with an abundance of 50% in the standard spectrum, the corresponding sample ion abundance must be between 30-70%).
c. Ions present at greater than 10% in the sample mass spectrum, but not present in the standard spectrum, must be evaluated by a reviewer experienced in mass spectral interpretation.

List compounds not meeting the criteria described above:

Sample ID	Compounds	Actions

Action:

1. The application of qualitative criteria for GC/MS analysis of target compounds requires professional judgment. It is up to the reviewer's discretion to obtain additional information from the laboratory. If it is determined that incorrect identifications were made, qualify all such data as unusable (R).

2. Use professional judgment to qualify the data if it is determined that cross-contamination has occurred.

3. Note in the Data Review Narrative any changes made to the reported compounds or concerns regarding target compound identifications. Note, for Contract Laboratory COR action, the necessity for numerous or significant changes.

TENTATIVELY IDENTIFIED COMPOUNDS (TICS)

NOTE: Tentatively identified compounds should only be evaluated when requested by a party from outside of the Hazardous Waste Support Section (HWSS).

List TICs

Sample ID	Compound	Sample ID	Compound

Action:

1. Qualify all TIC results for which there is presumptive evidence of a match (e.g. greater than or equal to 85% match) as tentatively identified (NJ), with approximated concentrations. TICs labeled "unknown" are qualified as estimated (J).
2. General actions related to the review of TIC results are as follows:
   a. If it is determined that a tentative identification of a non-target compound is unacceptable, change the tentative identification to "unknown" or another appropriate identification, and qualify the result as estimated (J).
   b. If all contractually-required peaks were not library searched and quantitated, the Region's designated representative may request these data from the laboratory.
3. In deciding whether a library search result for a TIC represents a reasonable identification, use professional judgment. If there is more than one possible match, report the result as "either compound X or compound Y". If there is a lack of isomer specificity, change the TIC result to a nonspecific isomer result (e.g., 1,3,5-trimethyl benzene to trimethyl benzene

isomer) or to a compound class (e.g., 2-methyl, 3-ethyl benzene to a substituted aromatic compound).

4. The reviewer may elect to report all similar compounds as a total (e.g., all alkanes may be summarized and reported as total hydrocarbons).
5. Target compounds from other fractions and suspected laboratory contaminants should be marked as "non-reportable".
6. Other Case factors may influence TIC judgments. If a sample TIC match is poor, but other samples have a TIC with a valid library match, similar RRT, and the same ions, infer identification information from the other sample TIC results.
7. Note in the Data Review Narrative any changes made to the reported data or any concerns regarding TIC identifications.
8. Note, for Contract Laboratory COR action, failure to properly evaluate and report TICs

All criteria were met _X_
Criteria were not met
and/or see below _____

SAMPLE QUANTITATION AND REPORTED CONTRACT REQUIRED QUANTITATION LIMITS (CRQLS)

Action:

1. If any discrepancies are found, the Region's designated representative may contact the laboratory to obtain additional information that could resolve any differences. If a discrepancy remains unresolved, the reviewer must use professional judgment to decide which value is the most accurate. Under these circumstances, the reviewer may determine that qualification of data is warranted. Note in the Data Review Narrative a description of the reasons for data qualification and the qualification that is applied to the data.
2. For non-aqueous samples, in the percent moisture is less than 70.0%, no qualification of the data is necessary. If the percent moisture is greater than or equal to 70.0% and less than 90.0%, qualify detects as estimated (J) and non-detects as approximated (UJ). If the percent moisture is greater than or equal to 90.0%, qualify detects as estimated (J) and non-detects as unusable (R) (see Table below).
3. Note, for Contract Laboratory COR action, numerous or significant failures to accurately quantify the target compounds or to properly evaluate and adjust CRQLs.
4. Results between MDL and CRQL should be qualified as estimated "J".
5. Results < MDL should be reported at the CRQL and qualified "U". MDLs themselves are not reported.

Table. Percent Moisture Actions for Low/Medium Volatiles Analysis for Non-Aqueous Samples

Criteria	**Action**	
	Detected Associated Compounds	**Non-detected Associated Compounds**
% Moisture < 70.0	No qualification	
70.0 < % Moisture < 90.0	J	UJ
% Moisture > 90.0	J	R

The sample quantitation evaluation is to verify laboratory quantitation results. In the space below, please show a minimum of one sample calculation:

Sample ID

FA42192-8 Isopropylbenzene RF = 1.728

[ ] = (1756934)(50)/(1.728)(1379039) = 36.86 ppb Ok

All criteria were met __X__
Criteria were not met
and/or see below _____

B. Percent Solids

List samples which have ≥ 70 % solids

_____
_____
_____
_____
_____
_____

QUANTITATION LIMITS

A. Dilution performed

SAMPLE ID	DILUTION FACTOR	REASON FOR DILUTION
FA42192-6	200 x; 2000 x	Several analytes above the calibration concentration range.

All criteria were met __X__
Criteria were not met
and/or see below _____

OTHER ISSUES

A. System Performance

List samples qualified based on the degradation of system performance during simple analysis:

Sample ID	Comments	Actions
_____	_____	_____
_No_degradation_of_system_performance_observed.		_____
_____	_____	_____
_____	_____	_____

Action:

Use professional judgment to qualify the data if it is determined that system performance has degraded during sample analyses. Inform the Contract Laboratory Program COR any action as a result of degradation of system performance which significantly affected the data.

B. Overall Assessment of Data

List samples qualified based on other issues:

Sample ID	Comments	Actions
_____	_____	_____
_No_additional_issues_observed_that_require_qualification_of_the_data._Results_are_valid_and__ _can_be_used_for_decission_purposes._____		_____
_____	_____	_____
_____	_____	_____

Action:

1. Use professional judgment to determine if there is any need to qualify data which were not qualified based on the Quality Control (QC) criteria previously discussed.
2. Write a brief narrative to give the user an indication of the analytical limitations of the data. Inform the Contract Laboratory COR the action, any inconsistency of the data with the Sample Delivery Group (SDG) Narrative. If sufficient information on the intended use and required quality of the data is available, the reviewer should include their assessment of the usability of the data within the given context. This may be used as part of a formal Data Quality Assessment (DQA).

EXECUTIVE NARRATIVE

SDG No:	FA42192	Laboratory:	Accutest, Orlando
Analysis:	SW846-8270D	Number of Samples:	11 Scan/22 SIM
Location:	BMSMC, Humacao, PR		

SUMMARY: Eleven (11) samples were analyzed for semivolatile organic compounds following method SW846-8270D. Twenty-two (22) samples were analyzed for selected PAHs and 1,4-Dioxane by SW846-8270D using the selective ion monitoring (SIM) technique; samples were analyzed separately for each analyte group. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence: EPA Hazardous Waste Support Section, SOP HW-35A, July 2015 – Revision 0. Semivolatile Data Validation. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

Results are valid and can be used for decision making purposes.

Critical issues: **None**
Major: **None**
Minor: **None**

Critical findings: **None**
Major findings: **None**
Minor findings:

1. All samples extracted and analyzed within method recommended holding time except for the cases described in the Data Review Worksheet. Sample preservation was appropriate.

Samples FA42192-12; FA42192-13; and FA42192-14 were re-extracted beyond holding time for confirmation run for surrogate recovery outside control limits. Results are not qualified, professional judgment.

2. Initial and continuing calibration verifications meet the method and guidance document required performance criteria except for the cases described in the Data Review Worksheet.

Analytes with % difference outside the method performance criteria but within guidance document performance criteria in the continuing calibration verification guidance document were not qualified.

No closing calibration verification included in data package for instrument GCMSL (SCAN). No action taken, professional judgment.

Analytes not meeting the initial and continuing % difference in the initial and continuing calibration verification criteria qualified as estimated (J or UJ) in affected samples.

3. No target analytes detected in method blanks except for the cases described in the Data Review Worksheet.

Laboratory qualified the results in samples FA42192-8, FA42192-11, FA42192-12, FA42192-13, FA42192-14, and FA42192-15 with a B qualifier.

Validator qualification based on guidelines: - Positive results below the reporting limits are reported at the reporting limits and qualified as non-detected (U). Results above the reporting limits and below method blanks results are reported at sample results and qualified as non-detects (U). Results above the reporting limits and blank result are not qualified, professional judgment.

4. DMCs meet the required criteria in all samples analyzed except for the cases described in the Data Review Worksheet. Non-deuterated surrogates added to the samples and were within laboratory recovery limits except in the cases described in the Data Review Worksheet;

Surrogate standards not recovered in sample FA42192-13 (SIM) due to dilution. No action taken, professional judgment. Surrogates recovered in the undiluted sample.

No qualification made on sample FA42192-6, professional judgment.

Terphenyl-d14 recovered low in samples FA42192-12 (31 %); FA42192-13 (32 %); and FA42192-14 (29 %); control limits 39 – 121 %. Results confirmed by re-extraction and analysis. Samples extracted beyond holding time. No action taken, professional judgment.

Surrogate recoveries in samples analyzed by SW846-8270D (SIM) were corrected for the concentration of the surrogate spiked. The correction is due to the following: the laboratory split batches between 8270 and 8270SIM. The surrogates were spiked at 8270 levels, which is 5x higher than what the laboratory 8270SIM setup expects. SIM surrogates recoveries were divided by 5. Calculations were spot checked.

5. As per chain-of-custody form sample FA42192-3 was to be used as the QC sample (MS/MSD) for this batch. Samples FA42192-6; FA42192-13; FA42157-7; and FA42237-15 were used instead. No action taken, professional judgment.

MS/MSD % recovery and RPD in sample FA42192-6 within laboratory control limits except for the cases described in the Data Review Worksheet.

Results for samples having either MS or the MSD outside the laboratory control limits are not qualified, professional judgment.

Samples with MS/MSD % recovery high are not qualified, professional judgment. 1,4-dioxane not qualified. No positive results associated with sample batch.

MS/MSD % recovery and RPD within laboratory control limits except for the cases described in the Data Review Worksheet.

Results for samples having either MS or the MSD outside the laboratory control limits are not qualified, professional judgment.

COMMENTS: Results are valid and can be used for decision making purposes.

Reviewers Name: Rafael Infante
Chemist License 1888

Signature: Rafael Infante

Date: April 26, 2017

SAMPLE ORGANIC DATA SAMPLE SUMMARY

Sample ID: FA42192-1
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: AQ - Field Blank Water

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	9.8	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	4.9	ug/l	1	-	U	Yes
2-Chlorophenol	4.9	ug/l	1	-	U	Yes
2,4-Dichlorophenol	4.9	ug/l	1	-	U	Yes
2,4-Dimethylphenol	4.9	ug/l	1	-	U	Yes
2,4-Dinitrophenol	25	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	9.8	ug/l	1	-	UJ	Yes
2-Methylphenol	4.9	ug/l	1	-	U	Yes
3&4-Methylphenol	4.9	ug/l	1	-	U	Yes
2-Nitrophenol	4.9	ug/l	1	-	UJ	Yes
4-Nitrophenol	25	ug/l	1	-	U	Yes
Pentachlorophenol	25	ug/l	1	-	U	Yes
Phenol	4.9	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	4.9	ug/l	1	-	UJ	Yes
2,4,5-Trichlorophenol	4.9	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	4.9	ug/l	1	-	U	Yes
Acenaphthene	4.9	ug/l	1	-	U	Yes
Acenaphthylene	4.9	ug/l	1	-	U	Yes
Acetophenone	4.9	ug/l	1	-	U	Yes
Anthracene	4.9	ug/l	1	-	U	Yes
Atrazine	4.9	ug/l	1	-	U	Yes
Benzaldehyde	25	ug/l	1	-	U	Yes
Benzo(a)anthracene	4.9	ug/l	1	-	U	Yes
Benzo(a)pyrene	4.9	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	4.9	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	4.9	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	4.9	ug/l	1	-	U	Yes
1,1'-Biphenyl	4.9	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	4.9	ug/l	1	-	U	Yes
Butyl benzyl phthalate	4.9	ug/l	1	-	U	Yes
Carbazole	4.9	ug/l	1	-	U	Yes
4-Chloroaniline	4.9	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	4.9	ug/l	1	-	U	Yes
bis(2-Chloroethyl)ether	4.9	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	4.9	ug/l	1	-	U	Yes

2-Chloronaphthalene	4.9	ug/l	1	-	U	Yes
4-Chlorophenyl phenyl ether	4.9	ug/l	1	-	U	Yes
Chrysene	4.9	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	4.9	ug/l	1	-	U	Yes
Dibenzofuran	4.9	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	4.9	ug/l	1	-	U	Yes
Diethyl phthalate	4.9	ug/l	1	-	U	Yes
Dimethyl phthalate	4.9	ug/l	1	-	U	Yes
Di-n-butyl phthalate	4.9	ug/l	1	-	U	Yes
Di-n-octyl phthalate	4.9	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	4.9	ug/l	1	-	UJ	Yes
2,6-Dinitrotoluene	4.9	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	4.9	ug/l	1	-	U	Yes
Fluoranthene	4.9	ug/l	1	-	U	Yes
Fluorene	4.9	ug/l	1	-	U	Yes
Hexachlorobenzene	4.9	ug/l	1	-	U	Yes
Hexachlorobutadiene	4.9	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	4.9	ug/l	1	-	U	Yes
Hexachloroethane	4.9	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	4.9	ug/l	1	-	U	Yes
Isophorone	4.9	ug/l	1	-	U	Yes
2-Methylnaphthalene	4.9	ug/l	1	-	U	Yes
2-Nitroaniline	4.9	ug/l	1	-	UJ	Yes
3-Nitroaniline	4.9	ug/l	1	-	U	Yes
4-Nitroaniline	4.9	ug/l	1	-	U	Yes
Nitrobenzene	4.9	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	4.9	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	4.9	ug/l	1	-	U	Yes
Phenanthrene	4.9	ug/l	1	-	U	Yes
Pyrene	4.9	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	4.9	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.20	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.20	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.20	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.20	ug/l	1	-	U	Yes
Chrysene	0.20	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.20	ug/l	1	-	U	Yes
1,4-Dioxane	0.29	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	0.20	ug/l	1	-	U	Yes
1-Methylnaphthalene	0.98	ug/l	1	-	U	Yes
2-Methylnaphthalene	0.98	ug/l	1	-	U	Yes
Naphthalene	0.98	ug/l	1	-	U	Yes

Sample ID: FA42192-2
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	10	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	5.0	ug/l	1	-	U	Yes
2-Chlorophenol	5.0	ug/l	1	-	U	Yes
2,4-Dichlorophenol	5.0	ug/l	1	-	U	Yes
2,4-Dimethylphenol	5.0	ug/l	1	-	U	Yes
2,4-Dinitrophenol	25	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	10	ug/l	1	-	UJ	Yes
2-Methylphenol	5.0	ug/l	1	-	U	Yes
3&4-Methylphenol	5.0	ug/l	1	-	U	Yes
2-Nitrophenol	5.0	ug/l	1	-	UJ	Yes
4-Nitrophenol	25	ug/l	1	-	U	Yes
Pentachlorophenol	25	ug/l	1	-	U	Yes
Phenol	5.0	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	5.0	ug/l	1	-	UJ	Yes
2,4,5-Trichlorophenol	5.0	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	5.0	ug/l	1	-	U	Yes
Acenaphthene	5.0	ug/l	1	-	U	Yes
Acenaphthylene	5.0	ug/l	1	-	U	Yes
Acetophenone	5.0	ug/l	1	-	U	Yes
Anthracene	1.3	ug/l	1	-	U	Yes
Atrazine	5.0	ug/l	1	-	U	Yes
Benzaldehyde	25	ug/l	1	-	U	Yes
Benzo(a)anthracene	5.0	ug/l	1	-	U	Yes
Benzo(a)pyrene	5.0	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	5.0	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	5.0	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	5.0	ug/l	1	-	U	Yes
1,1'-Biphenyl	5.0	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	5.0	ug/l	1	-	U	Yes
Butyl benzyl phthalate	5.0	ug/l	1	-	U	Yes
Carbazole	5.0	ug/l	1	-	U	Yes
4-Chloroaniline	5.0	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	5.0	ug/l	1	-	U	Yes
bis(2-Chloroethyl)ether	5.0	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	5.0	ug/l	1	-	U	Yes
2-Chloronaphthalene	5.0	ug/l	1	-	U	Yes
4-Chlorophenyl phenyl ether	5.0	ug/l	1	-	U	Yes

Chrysene	5.0	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	5.0	ug/l	1	-	U	Yes
Dibenzofuran	5.0	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	5.0	ug/l	1	-	U	Yes
Diethyl phthalate	5.0	ug/l	1	-	U	Yes
Dimethyl phthalate	5.0	ug/l	1	-	U	Yes
Di-n-butyl phthalate	5.0	ug/l	1	-	U	Yes
Di-n-octyl phthalate	5.0	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	5.0	ug/l	1	-	UJ	Yes
2,6-Dinitrotoluene	5.0	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	5.0	ug/l	1	-	U	Yes
Fluoranthene	5.0	ug/l	1	-	U	Yes
Fluorene	5.0	ug/l	1	-	U	Yes
Hexachlorobenzene	5.0	ug/l	1	-	U	Yes
Hexachlorobutadiene	5.0	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	5.0	ug/l	1	-	U	Yes
Hexachloroethane	5.0	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	5.0	ug/l	1	-	U	Yes
Isophorone	5.0	ug/l	1	-	U	Yes
2-Methylnaphthalene	5.0	ug/l	1	-	U	Yes
2-Nitroaniline	5.0	ug/l	1	-	UJ	Yes
3-Nitroaniline	5.0	ug/l	1	-	U	Yes
4-Nitroaniline	5.0	ug/l	1	-	U	Yes
Nitrobenzene	5.0	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	5.0	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	5.0	ug/l	1	-	U	Yes
Phenanthrene	5.0	ug/l	1	-	U	Yes
Pyrene	5.0	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	5.0	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.20	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.20	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.20	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.20	ug/l	1	-	U	Yes
Chrysene	0.20	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.20	ug/l	1	-	U	Yes
1,4-Dioxane	15.9	ug/l	1	-	-	Yes
Indeno(1,2,3-cd)pyrene	0.20	ug/l	1	-	U	Yes
1-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
2-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
Naphthalene	1.0	ug/l	1	-	U	Yes

Sample ID: FA42192-3
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	9.8	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	4.9	ug/l	1	-	U	Yes
2-Chlorophenol	4.9	ug/l	1	-	U	Yes
2,4-Dichlorophenol	4.9	ug/l	1	-	U	Yes
2,4-Dimethylphenol	4.9	ug/l	1	-	U	Yes
2,4-Dinitrophenol	25	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	9.8	ug/l	1	-	UJ	Yes
2-Methylphenol	4.9	ug/l	1	-	U	Yes
3&4-Methylphenol	4.9	ug/l	1	-	U	Yes
2-Nitrophenol	4.9	ug/l	1	-	UJ	Yes
4-Nitrophenol	25	ug/l	1	-	U	Yes
Pentachlorophenol	25	ug/l	1	-	U	Yes
Phenol	4.9	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	4.9	ug/l	1	-	UJ	Yes
2,4,5-Trichlorophenol	4.9	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	4.9	ug/l	1	-	U	Yes
Acenaphthene	4.9	ug/l	1	-	U	Yes
Acenaphthylene	4.9	ug/l	1	-	U	Yes
Acetophenone	4.9	ug/l	1	-	U	Yes
Anthracene	4.9	ug/l	1	-	U	Yes
Atrazine	4.9	ug/l	1	-	U	Yes
Benzaldehyde	25	ug/l	1	-	U	Yes
Benzo(a)anthracene	4.9	ug/l	1	-	U	Yes
Benzo(a)pyrene	4.9	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	4.9	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	4.9	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	4.9	ug/l	1	-	U	Yes
1,1'-Biphenyl	4.9	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	4.9	ug/l	1	-	U	Yes
Butyl benzyl phthalate	4.9	ug/l	1	-	U	Yes
Carbazole	4.9	ug/l	1	-	U	Yes
4-Chloroaniline	4.9	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	4.9	ug/l	1	-	U	Yes
bis(2-Chloroethyl)ether	4.9	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	4.9	ug/l	1	-	U	Yes
2-Chloronaphthalene	4.9	ug/l	1	-	U	Yes
4-Chlorophenyl phenyl ether	4.9	ug/l	1	-	U	Yes

Chrysene	4.9	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	4.9	ug/l	1	-	U	Yes
Dibenzofuran	4.9	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	4.9	ug/l	1	-	U	Yes
Diethyl phthalate	4.9	ug/l	1	-	U	Yes
Dimethyl phthalate	4.9	ug/l	1	-	U	Yes
Di-n-butyl phthalate	4.9	ug/l	1	-	U	Yes
Di-n-octyl phthalate	4.9	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	4.9	ug/l	1	-	UJ	Yes
2,6-Dinitrotoluene	4.9	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	4.9	ug/l	1	-	U	Yes
Fluoranthene	4.9	ug/l	1	-	U	Yes
Fluorene	4.9	ug/l	1	-	U	Yes
Hexachlorobenzene	4.9	ug/l	1	-	U	Yes
Hexachlorobutadiene	4.9	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	4.9	ug/l	1	-	U	Yes
Hexachloroethane	4.9	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	4.9	ug/l	1	-	U	Yes
Isophorone	4.9	ug/l	1	-	U	Yes
2-Methylnaphthalene	4.9	ug/l	1	-	U	Yes
2-Nitroaniline	4.9	ug/l	1	-	UJ	Yes
3-Nitroaniline	4.9	ug/l	1	-	U	Yes
4-Nitroaniline	4.9	ug/l	1	-	U	Yes
Nitrobenzene	4.9	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	4.9	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	4.9	ug/l	1	-	U	Yes
Phenanthrene	1.5	ug/l	1	J	J	Yes
Pyrene	4.9	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	4.9	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.20	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.20	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.20	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.20	ug/l	1	-	U	Yes
Chrysene	0.20	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.20	ug/l	1	-	U	Yes
1,4-Dioxane	18.4	ug/l	1	-	-	Yes
Indeno(1,2,3-cd)pyrene	0.20	ug/l	1	-	U	Yes
1-Methylnaphthalene	0.98	ug/l	1	-	U	Yes
2-Methylnaphthalene	0.98	ug/l	1	-	U	Yes
Naphthalene	0.98	ug/l	1	-	U	Yes

Sample ID: FA42192-6
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	100	ug/l	10	-	U	Yes
4-Chloro-3-methyl phenol	50	ug/l	10	-	U	Yes
2-Chlorophenol	50	ug/l	10	-	U	Yes
2,4-Dichlorophenol	50	ug/l	10	-	U	Yes
2,4-Dimethylphenol	50	ug/l	10	-	U	Yes
2,4-Dinitrophenol	250	ug/l	10	-	U	Yes
4,6-Dinitro-o-cresol	100	ug/l	10	-	UJ	Yes
2-Methylphenol	50	ug/l	10	-	U	Yes
3&4-Methylphenol	50	ug/l	10	-	U	Yes
2-Nitrophenol	50	ug/l	10	-	UJ	Yes
4-Nitrophenol	250	ug/l	10	-	U	Yes
Pentachlorophenol	250	ug/l	10	-	U	Yes
Phenol	50	ug/l	10	-	U	Yes
2,3,4,6-Tetrachlorophenol	50	ug/l	10	-	UJ	Yes
2,4,5-Trichlorophenol	50	ug/l	10	-	U	Yes
2,4,6-Trichlorophenol	50	ug/l	10	-	U	Yes
Acenaphthene	50	ug/l	10	-	U	Yes
Acenaphthylene	50	ug/l	10	-	U	Yes
Acetophenone	50	ug/l	10	-	U	Yes
Anthracene	27.9	ug/l	10	J	J	Yes
Atrazine	50	ug/l	10	-	U	Yes
Benzaldehyde	250	ug/l	10	-	U	Yes
Benzo(a)anthracene	50	ug/l	10	-	U	Yes
Benzo(a)pyrene	50	ug/l	10	-	U	Yes
Benzo(b)fluoranthene	50	ug/l	10	-	U	Yes
Benzo(g,h,i)perylene	50	ug/l	10	-	U	Yes
Benzo(k)fluoranthene	50	ug/l	10	-	U	Yes
1,1'-Biphenyl	50	ug/l	10	-	U	Yes
4-Bromophenyl phenyl ether	50	ug/l	10	-	U	Yes
Butyl benzyl phthalate	50	ug/l	10	-	U	Yes
Carbazole	50	ug/l	10	-	U	Yes
4-Chloroaniline	50	ug/l	10	-	U	Yes
bis(2-Chloroethoxy)methane	50	ug/l	10	-	U	Yes
bis(2-Chloroethyl)ether	50	ug/l	10	-	U	Yes
bis(2-Chloroisopropyl)ether	50	ug/l	10	-	U	Yes
2-Chloronaphthalene	50	ug/l	10	-	U	Yes
4-Chlorophenyl phenyl ether	50	ug/l	10	-	U	Yes

Chrysene	50	ug/l	10	-	U	Yes
Dibenzo(a,h)anthracene	50	ug/l	10	-	U	Yes
Dibenzofuran	50	ug/l	10	-	U	Yes
3,3'-Dichlorobenzidine	50	ug/l	10	-	U	Yes
Diethyl phthalate	50	ug/l	10	-	U	Yes
Dimethyl phthalate	50	ug/l	10	-	U	Yes
Di-n-butyl phthalate	50	ug/l	10	-	U	Yes
Di-n-octyl phthalate	50	ug/l	10	-	U	Yes
2,4-Dinitrotoluene	50	ug/l	10	-	UJ	Yes
2,6-Dinitrotoluene	50	ug/l	10	-	U	Yes
bis(2-Ethylhexyl)phthalate	50	ug/l	10	-	U	Yes
Fluoranthene	50	ug/l	10	-	U	Yes
Fluorene	50	ug/l	10	-	U	Yes
Hexachlorobenzene	50	ug/l	10	-	U	Yes
Hexachlorobutadiene	50	ug/l	10	-	U	Yes
Hexachlorocyclopentadiene	50	ug/l	10	-	U	Yes
Hexachloroethane	50	ug/l	10	-	U	Yes
Indeno(1,2,3-cd)pyrene	50	ug/l	10	-	U	Yes
Isophorone	50	ug/l	10	-	U	Yes
2-Methylnaphthalene	50	ug/l	10	-	U	Yes
2-Nitroaniline	50	ug/l	10	-	UJ	Yes
3-Nitroaniline	50	ug/l	10	-	U	Yes
4-Nitroaniline	50	ug/l	10	-	U	Yes
Nitrobenzene	50	ug/l	10	-	U	Yes
N-Nitroso-di-n-propylamine	50	ug/l	10	-	U	Yes
N-Nitrosodiphenylamine	50	ug/l	10	-	U	Yes
Phenanthrene	50	ug/l	10	-	U	Yes
Pyrene	50	ug/l	10	-	U	Yes
1,2,4,5-Tetrachlorobenzene	50	ug/l	10	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.20	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.20	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.20	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.20	ug/l	1	-	U	Yes
Chrysene	0.20	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.20	ug/l	1	-	U	Yes
1,4-Dioxane	3.0	ug/l	1	-	-	Yes
Indeno(1,2,3-cd)pyrene	0.20	ug/l	1	-	U	Yes
1-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
2-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
Naphthalene	1.0	ug/l	1	-	U	Yes

Sample ID: FA42192-7
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	9.6	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	4.8	ug/l	1	-	U	Yes
2-Chlorophenol	4.8	ug/l	1	-	U	Yes
2,4-Dichlorophenol	4.8	ug/l	1	-	U	Yes
2,4-Dimethylphenol	4.8	ug/l	1	-	U	Yes
2,4-Dinitrophenol	24	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	9.6	ug/l	1	-	UJ	Yes
2-Methylphenol	4.8	ug/l	1	-	U	Yes
3&4-Methylphenol	4.8	ug/l	1	-	U	Yes
2-Nitrophenol	4.8	ug/l	1	-	UJ	Yes
4-Nitrophenol	24	ug/l	1	-	U	Yes
Pentachlorophenol	24	ug/l	1	-	U	Yes
Phenol	4.8	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	4.8	ug/l	1	-	UJ	Yes
2,4,5-Trichlorophenol	4.8	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	4.8	ug/l	1	-	U	Yes
Acenaphthene	4.8	ug/l	1	-	U	Yes
Acenaphthylene	4.8	ug/l	1	-	U	Yes
Acetophenone	4.8	ug/l	1	-	U	Yes
Anthracene	1.3	ug/l	1	-	U	Yes
Atrazine	4.8	ug/l	1	-	U	Yes
Benzaldehyde	24	ug/l	1	-	U	Yes
Benzo(a)anthracene	4.8	ug/l	1	-	U	Yes
Benzo(a)pyrene	4.8	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	4.8	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	4.8	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	4.8	ug/l	1	-	U	Yes
1,1'-Biphenyl	4.8	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	4.8	ug/l	1	-	U	Yes
Butyl benzyl phthalate	4.8	ug/l	1	-	U	Yes
Carbazole	4.8	ug/l	1	-	U	Yes
4-Chloroaniline	4.8	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	4.8	ug/l	1	-	U	Yes
bis(2-Chloroethyl)ether	4.8	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	4.8	ug/l	1	-	U	Yes
2-Chloronaphthalene	4.8	ug/l	1	-	U	Yes

4-Chlorophenyl phenyl ether	4.8	ug/l	1	-	U	Yes
Chrysene	4.8	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	4.8	ug/l	1	-	U	Yes
Dibenzofuran	4.8	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	4.8	ug/l	1	-	U	Yes
Diethyl phthalate	4.8	ug/l	1	-	U	Yes
Dimethyl phthalate	4.8	ug/l	1	-	U	Yes
Di-n-butyl phthalate	4.8	ug/l	1	-	U	Yes
Di-n-octyl phthalate	4.8	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	4.8	ug/l	1	-	UJ	Yes
2,6-Dinitrotoluene	4.8	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	4.8	ug/l	1	-	U	Yes
Fluoranthene	4.8	ug/l	1	-	U	Yes
Fluorene	4.8	ug/l	1	-	U	Yes
Hexachlorobenzene	4.8	ug/l	1	-	U	Yes
Hexachlorobutadiene	4.8	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	4.8	ug/l	1	-	U	Yes
Hexachloroethane	4.8	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	4.8	ug/l	1	-	U	Yes
Isophorone	4.8	ug/l	1	-	U	Yes
2-Methylnaphthalene	4.8	ug/l	1	-	U	Yes
2-Nitroaniline	4.8	ug/l	1	-	UJ	Yes
3-Nitroaniline	4.8	ug/l	1	-	U	Yes
4-Nitroaniline	4.8	ug/l	1	-	U	Yes
Nitrobenzene	4.8	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	4.8	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	4.8	ug/l	1	-	U	Yes
Phenanthrene	4.8	ug/l	1	-	U	Yes
Pyrene	4.8	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	4.8	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.19	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.19	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.19	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.19	ug/l	1	-	U	Yes
Chrysene	0.19	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.19	ug/l	1	-	U	Yes
1,4-Dioxane	2.6	ug/l	1	-	-	Yes
Indeno(1,2,3-cd)pyrene	0.19	ug/l	1	-	U	Yes
1-Methylnaphthalene	0.96	ug/l	1	-	U	Yes
2-Methylnaphthalene	0.96	ug/l	1	-	U	Yes
Naphthalene	0.96	ug/l	1	-	U	Yes

Sample ID: FA42192-8
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	10	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	5.0	ug/l	1	-	U	Yes
2-Chlorophenol	5.0	ug/l	1	-	U	Yes
2,4-Dichlorophenol	5.0	ug/l	1	-	U	Yes
2,4-Dimethylphenol	5.0	ug/l	1	-	U	Yes
2,4-Dinitrophenol	25	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	10	ug/l	1	-	U	Yes
2-Methylphenol	5.0	ug/l	1	-	U	Yes
3&4-Methylphenol	5.0	ug/l	1	-	U	Yes
2-Nitrophenol	5.0	ug/l	1	-	U	Yes
4-Nitrophenol	25	ug/l	1	-	U	Yes
Pentachlorophenol	25	ug/l	1	-	U	Yes
Phenol	5.0	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	5.0	ug/l	1	-	U	Yes
2,4,5-Trichlorophenol	5.0	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	5.0	ug/l	1	-	U	Yes
Acenaphthene	5.0	ug/l	1	-	U	Yes
Acenaphthylene	5.0	ug/l	1	-	U	Yes
Acetophenone	5.0	ug/l	1	-	U	Yes
Anthracene	1.3	ug/l	1	J	J	Yes
Atrazine	5.0	ug/l	1	-	U	Yes
Benzaldehyde	25	ug/l	1	-	U	Yes
Benzo(a)anthracene	5.0	ug/l	1	-	U	Yes
Benzo(a)pyrene	5.0	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	5.0	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	5.0	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	5.0	ug/l	1	-	U	Yes
1,1'-Biphenyl	5.0	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	5.0	ug/l	1	-	U	Yes
Butyl benzyl phthalate	5.0	ug/l	1	-	U	Yes
Carbazole	5.0	ug/l	1	-	U	Yes
4-Chloroaniline	5.0	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	5.0	ug/l	1	-	UJ	Yes
bis(2-Chloroethyl)ether	5.0	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	5.0	ug/l	1	-	U	Yes
2-Chloronaphthalene	5.0	ug/l	1	-	U	Yes

4-Chlorophenyl phenyl ether	5.0	ug/l	1	-	U	Yes
Chrysene	5.0	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	5.0	ug/l	1	-	U	Yes
Dibenzofuran	5.0	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	5.0	ug/l	1	-	U	Yes
Diethyl phthalate	5.0	ug/l	1	-	U	Yes
Dimethyl phthalate	5.0	ug/l	1	-	U	Yes
Di-n-butyl phthalate	5.0	ug/l	1	-	U	Yes
Di-n-octyl phthalate	5.0	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	5.0	ug/l	1	-	U	Yes
2,6-Dinitrotoluene	5.0	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	5.0	ug/l	1	-	U	Yes
Fluoranthene	5.0	ug/l	1	-	U	Yes
Fluorene	5.0	ug/l	1	-	U	Yes
Hexachlorobenzene	5.0	ug/l	1	-	U	Yes
Hexachlorobutadiene	5.0	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	5.0	ug/l	1	-	U	Yes
Hexachloroethane	5.0	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	5.0	ug/l	1	-	U	Yes
Isophorone	5.0	ug/l	1	-	U	Yes
2-Methylnaphthalene	5.0	ug/l	1	-	UJ	Yes
2-Nitroaniline	5.0	ug/l	1	-	U	Yes
3-Nitroaniline	5.0	ug/l	1	-	U	Yes
4-Nitroaniline	5.0	ug/l	1	-	U	Yes
Nitrobenzene	5.0	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	5.0	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	5.0	ug/l	1	-	U	Yes
Phenanthrene	5.0	ug/l	1	-	U	Yes
Pyrene	5.0	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	5.0	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.20	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.20	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.20	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.20	ug/l	1	-	U	Yes
Chrysene	0.20	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.20	ug/l	1	-	U	Yes
1,4-Dioxane	1.2	ug/l	1	B	-	Yes
Indeno(1,2,3-cd)pyrene	0.20	ug/l	1	-	U	Yes
1-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
2-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
Naphthalene	1.0	ug/l	1	-	U	Yes

Sample ID: FA42192-11
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	9.8	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	4.9	ug/l	1	-	U	Yes
2-Chlorophenol	4.9	ug/l	1	-	U	Yes
2,4-Dichlorophenol	4.9	ug/l	1	-	U	Yes
2,4-Dimethylphenol	4.9	ug/l	1	-	U	Yes
2,4-Dinitrophenol	25	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	9.8	ug/l	1	-	U	Yes
2-Methylphenol	4.9	ug/l	1	-	U	Yes
3&4-Methylphenol	4.9	ug/l	1	-	U	Yes
2-Nitrophenol	4.9	ug/l	1	-	U	Yes
4-Nitrophenol	25	ug/l	1	-	U	Yes
Pentachlorophenol	25	ug/l	1	-	U	Yes
Phenol	4.9	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	4.9	ug/l	1	-	U	Yes
2,4,5-Trichlorophenol	4.9	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	4.9	ug/l	1	-	U	Yes
Acenaphthene	4.9	ug/l	1	-	U	Yes
Acenaphthylene	4.9	ug/l	1	-	U	Yes
Acetophenone	4.9	ug/l	1	-	U	Yes
Anthracene	1.4	ug/l	1	J	J	Yes
Atrazine	4.9	ug/l	1	-	U	Yes
Benzaldehyde	25	ug/l	1	-	U	Yes
Benzo(a)anthracene	4.9	ug/l	1	-	U	Yes
Benzo(a)pyrene	4.9	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	4.9	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	4.9	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	4.9	ug/l	1	-	U	Yes
1,1'-Biphenyl	4.9	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	4.9	ug/l	1	-	U	Yes
Butyl benzyl phthalate	4.9	ug/l	1	-	U	Yes
Carbazole	4.9	ug/l	1	-	U	Yes
4-Chloroaniline	4.9	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	4.9	ug/l	1	-	UJ	Yes
bis(2-Chloroethyl)ether	4.9	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	4.9	ug/l	1	-	U	Yes
2-Chloronaphthalene	4.9	ug/l	1	-	U	Yes

4-Chlorophenyl phenyl ether	4.9	ug/l	1	-	U	Yes
Chrysene	4.9	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	4.9	ug/l	1	-	U	Yes
Dibenzofuran	4.9	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	4.9	ug/l	1	-	U	Yes
Diethyl phthalate	4.9	ug/l	1	-	U	Yes
Dimethyl phthalate	4.9	ug/l	1	-	U	Yes
Di-n-butyl phthalate	4.9	ug/l	1	-	U	Yes
Di-n-octyl phthalate	4.9	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	4.9	ug/l	1	-	U	Yes
2,6-Dinitrotoluene	4.9	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	4.9	ug/l	1	-	U	Yes
Fluoranthene	4.9	ug/l	1	-	U	Yes
Fluorene	4.9	ug/l	1	-	U	Yes
Hexachlorobenzene	4.9	ug/l	1	-	U	Yes
Hexachlorobutadiene	4.9	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	4.9	ug/l	1	-	U	Yes
Hexachloroethane	4.9	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	4.9	ug/l	1	-	U	Yes
Isophorone	4.9	ug/l	1	-	U	Yes
2-Methylnaphthalene	4.9	ug/l	1	-	UJ	Yes
2-Nitroaniline	4.9	ug/l	1	-	U	Yes
3-Nitroaniline	4.9	ug/l	1	-	U	Yes
4-Nitroaniline	4.9	ug/l	1	-	U	Yes
Nitrobenzene	4.9	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	4.9	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	4.9	ug/l	1	-	U	Yes
Phenanthrene	4.9	ug/l	1	-	U	Yes
Pyrene	4.9	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	4.9	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.20	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.20	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.20	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.20	ug/l	1	-	U	Yes
Chrysene	0.20	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.20	ug/l	1	-	U	Yes
1,4-Dioxane	1.5	ug/l	1	B	-	Yes
Indeno(1,2,3-cd)pyrene	0.20	ug/l	1	-	U	Yes
1-Methylnaphthalene	0.98	ug/l	1	-	U	Yes
2-Methylnaphthalene	0.98	ug/l	1	-	U	Yes
Naphthalene	0.98	ug/l	1	-	U	Yes

Sample ID: FA42192-12
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	9.8	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	4.9	ug/l	1	-	U	Yes
2-Chlorophenol	4.9	ug/l	1	-	U	Yes
2,4-Dichlorophenol	4.9	ug/l	1	-	U	Yes
2,4-Dimethylphenol	4.9	ug/l	1	-	U	Yes
2,4-Dinitrophenol	25	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	9.8	ug/l	1	-	U	Yes
2-Methylphenol	4.9	ug/l	1	-	U	Yes
3&4-Methylphenol	4.9	ug/l	1	-	U	Yes
2-Nitrophenol	4.9	ug/l	1	-	U	Yes
4-Nitrophenol	25	ug/l	1	-	U	Yes
Pentachlorophenol	25	ug/l	1	-	U	Yes
Phenol	4.9	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	4.9	ug/l	1	-	U	Yes
2,4,5-Trichlorophenol	4.9	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	4.9	ug/l	1	-	U	Yes
Acenaphthene	4.9	ug/l	1	-	U	Yes
Acenaphthylene	4.9	ug/l	1	-	U	Yes
Acetophenone	4.9	ug/l	1	-	U	Yes
Anthracene	4.9	ug/l	1	-	U	Yes
Atrazine	4.9	ug/l	1	-	U	Yes
Benzaldehyde	25	ug/l	1	-	U	Yes
Benzo(a)anthracene	4.9	ug/l	1	-	U	Yes
Benzo(a)pyrene	4.9	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	4.9	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	4.9	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	4.9	ug/l	1	-	U	Yes
1,1'-Biphenyl	4.9	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	4.9	ug/l	1	-	U	Yes
Butyl benzyl phthalate	4.9	ug/l	1	-	U	Yes
Carbazole	4.9	ug/l	1	-	U	Yes
4-Chloroaniline	4.9	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	4.9	ug/l	1	-	UJ	Yes
bis(2-Chloroethyl)ether	4.9	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	4.9	ug/l	1	-	U	Yes
2-Chloronaphthalene	4.9	ug/l	1	-	U	Yes

4-Chlorophenyl phenyl ether	4.9	ug/l	1	-	U	Yes
Chrysene	4.9	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	4.9	ug/l	1	-	U	Yes
Dibenzofuran	4.9	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	4.9	ug/l	1	-	U	Yes
Diethyl phthalate	4.9	ug/l	1	-	U	Yes
Dimethyl phthalate	4.9	ug/l	1	-	U	Yes
Di-n-butyl phthalate	4.9	ug/l	1	-	U	Yes
Di-n-octyl phthalate	4.9	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	4.9	ug/l	1	-	U	Yes
2,6-Dinitrotoluene	4.9	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	4.9	ug/l	1	-	U	Yes
Fluoranthene	4.9	ug/l	1	-	U	Yes
Fluorene	4.9	ug/l	1	-	U	Yes
Hexachlorobenzene	4.9	ug/l	1	-	U	Yes
Hexachlorobutadiene	4.9	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	4.9	ug/l	1	-	U	Yes
Hexachloroethane	4.9	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	4.9	ug/l	1	-	U	Yes
Isophorone	4.9	ug/l	1	-	U	Yes
2-Methylnaphthalene	4.9	ug/l	1	-	UJ	Yes
2-Nitroaniline	4.9	ug/l	1	-	U	Yes
3-Nitroaniline	4.9	ug/l	1	-	U	Yes
4-Nitroaniline	4.9	ug/l	1	-	U	Yes
Nitrobenzene	4.9	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	4.9	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	4.9	ug/l	1	-	U	Yes
Phenanthrene	4.9	ug/l	1	-	U	Yes
Pyrene	4.9	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	4.9	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.20	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.20	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.20	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.20	ug/l	1	-	U	Yes
Chrysene	0.20	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.20	ug/l	1	-	U	Yes
1,4-Dioxane	103	ug/l	10	B	-	Yes
Indeno(1,2,3-cd)pyrene	0.20	ug/l	1	-	U	Yes
1-Methylnaphthalene	0.98	ug/l	1	-	U	Yes
2-Methylnaphthalene	0.98	ug/l	1	-	U	Yes
Naphthalene	0.98	ug/l	1	-	U	Yes

Sample ID: FA42192-13
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	10	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	5.0	ug/l	1	-	U	Yes
2-Chlorophenol	5.0	ug/l	1	-	U	Yes
2,4-Dichlorophenol	5.0	ug/l	1	-	U	Yes
2,4-Dimethylphenol	5.0	ug/l	1	-	U	Yes
2,4-Dinitrophenol	25	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	10	ug/l	1	-	U	Yes
2-Methylphenol	5.0	ug/l	1	-	U	Yes
3&4-Methylphenol	5.0	ug/l	1	-	U	Yes
2-Nitrophenol	5.0	ug/l	1	-	U	Yes
4-Nitrophenol	25	ug/l	1	-	U	Yes
Pentachlorophenol	25	ug/l	1	-	U	Yes
Phenol	5.0	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	5.0	ug/l	1	-	U	Yes
2,4,5-Trichlorophenol	5.0	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	5.0	ug/l	1	-	U	Yes
Acenaphthene	5.0	ug/l	1	-	U	Yes
Acenaphthylene	5.0	ug/l	1	-	U	Yes
Acetophenone	5.0	ug/l	1	-	U	Yes
Anthracene	5.0	ug/l	1	-	U	Yes
Atrazine	5.0	ug/l	1	-	U	Yes
Benzaldehyde	25	ug/l	1	-	U	Yes
Benzo(a)anthracene	5.0	ug/l	1	-	U	Yes
Benzo(a)pyrene	5.0	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	5.0	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	5.0	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	5.0	ug/l	1	-	U	Yes
1,1'-Biphenyl	5.0	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	5.0	ug/l	1	-	U	Yes
Butyl benzyl phthalate	5.0	ug/l	1	-	U	Yes
Carbazole	5.0	ug/l	1	-	U	Yes
4-Chloroaniline	5.0	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	5.0	ug/l	1	-	UJ	Yes
bis(2-Chloroethyl)ether	5.0	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	5.0	ug/l	1	-	U	Yes
2-Chloronaphthalene	5.0	ug/l	1	-	U	Yes

4-Chlorophenyl phenyl ether	5.0	ug/l	1	-	U	Yes
Chrysene	5.0	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	5.0	ug/l	1	-	U	Yes
Dibenzofuran	5.0	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	5.0	ug/l	1	-	U	Yes
Diethyl phthalate	5.0	ug/l	1	-	U	Yes
Dimethyl phthalate	5.0	ug/l	1	-	U	Yes
Di-n-butyl phthalate	5.0	ug/l	1	-	U	Yes
Di-n-octyl phthalate	5.0	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	5.0	ug/l	1	-	U	Yes
2,6-Dinitrotoluene	5.0	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	5.0	ug/l	1	-	U	Yes
Fluoranthene	5.0	ug/l	1	-	U	Yes
Fluorene	5.0	ug/l	1	-	U	Yes
Hexachlorobenzene	5.0	ug/l	1	-	U	Yes
Hexachlorobutadiene	5.0	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	5.0	ug/l	1	-	U	Yes
Hexachloroethane	5.0	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	5.0	ug/l	1	-	U	Yes
Isophorone	5.0	ug/l	1	-	U	Yes
2-Methylnaphthalene	5.0	ug/l	1	-	UJ	Yes
2-Nitroaniline	5.0	ug/l	1	-	U	Yes
3-Nitroaniline	5.0	ug/l	1	-	U	Yes
4-Nitroaniline	5.0	ug/l	1	-	U	Yes
Nitrobenzene	5.0	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	5.0	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	5.0	ug/l	1	-	U	Yes
Phenanthrene	5.0	ug/l	1	-	U	Yes
Pyrene	5.0	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	5.0	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.20	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.20	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.20	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.20	ug/l	1	-	U	Yes
Chrysene	0.20	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.20	ug/l	1	-	U	Yes
1,4-Dioxane	313	ug/l	25	B	-	Yes
Indeno(1,2,3-cd)pyrene	0.20	ug/l	1	-	U	Yes
1-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
2-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
Naphthalene	1.0	ug/l	1	-	U	Yes

Sample ID: FA42192-14
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	9.6	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	4.8	ug/l	1	-	U	Yes
2-Chlorophenol	4.8	ug/l	1	-	U	Yes
2,4-Dichlorophenol	4.8	ug/l	1	-	U	Yes
2,4-Dimethylphenol	4.8	ug/l	1	-	U	Yes
2,4-Dinitrophenol	24	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	9.6	ug/l	1	-	U	Yes
2-Methylphenol	4.8	ug/l	1	-	U	Yes
3&4-Methylphenol	4.8	ug/l	1	-	U	Yes
2-Nitrophenol	4.8	ug/l	1	-	U	Yes
4-Nitrophenol	24	ug/l	1	-	U	Yes
Pentachlorophenol	24	ug/l	1	-	U	Yes
Phenol	4.8	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	4.8	ug/l	1	-	U	Yes
2,4,5-Trichlorophenol	4.8	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	4.8	ug/l	1	-	U	Yes
Acenaphthene	4.8	ug/l	1	-	U	Yes
Acenaphthylene	4.8	ug/l	1	-	U	Yes
Acetophenone	4.8	ug/l	1	-	U	Yes
Anthracene	1.3	ug/l	1	-	U	Yes
Atrazine	4.8	ug/l	1	-	U	Yes
Benzaldehyde	24	ug/l	1	-	U	Yes
Benzo(a)anthracene	4.8	ug/l	1	-	U	Yes
Benzo(a)pyrene	4.8	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	4.8	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	4.8	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	4.8	ug/l	1	-	U	Yes
1,1'-Biphenyl	4.8	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	4.8	ug/l	1	-	U	Yes
Butyl benzyl phthalate	4.8	ug/l	1	-	U	Yes
Carbazole	4.8	ug/l	1	-	U	Yes
4-Chloroaniline	4.8	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	4.8	ug/l	1	-	UJ	Yes
bis(2-Chloroethyl)ether	4.8	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	4.8	ug/l	1	-	U	Yes
2-Chloronaphthalene	4.8	ug/l	1	-	U	Yes

4-Chlorophenyl phenyl ether	4.8	ug/l	1	-	U	Yes
Chrysene	4.8	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	4.8	ug/l	1	-	U	Yes
Dibenzofuran	4.8	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	4.8	ug/l	1	-	U	Yes
Diethyl phthalate	4.8	ug/l	1	-	U	Yes
Dimethyl phthalate	4.8	ug/l	1	-	U	Yes
Di-n-butyl phthalate	4.8	ug/l	1	-	U	Yes
Di-n-octyl phthalate	4.8	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	4.8	ug/l	1	-	U	Yes
2,6-Dinitrotoluene	4.8	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	4.8	ug/l	1	-	U	Yes
Fluoranthene	4.8	ug/l	1	-	U	Yes
Fluorene	4.8	ug/l	1	-	U	Yes
Hexachlorobenzene	4.8	ug/l	1	-	U	Yes
Hexachlorobutadiene	4.8	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	4.8	ug/l	1	-	U	Yes
Hexachloroethane	4.8	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	4.8	ug/l	1	-	U	Yes
Isophorone	4.8	ug/l	1	-	U	Yes
2-Methylnaphthalene	4.8	ug/l	1	-	UJ	Yes
2-Nitroaniline	4.8	ug/l	1	-	U	Yes
3-Nitroaniline	4.8	ug/l	1	-	U	Yes
4-Nitroaniline	4.8	ug/l	1	-	U	Yes
Nitrobenzene	4.8	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	4.8	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	4.8	ug/l	1	-	U	Yes
Phenanthrene	4.8	ug/l	1	-	U	Yes
Pyrene	4.8	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	4.8	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.19	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.19	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.19	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.19	ug/l	1	-	U	Yes
Chrysene	0.19	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.19	ug/l	1	-	U	Yes
1,4-Dioxane	15.8	ug/l	1	B	-	Yes
Indeno(1,2,3-cd)pyrene	0.19	ug/l	1	-	U	Yes
1-Methylnaphthalene	0.96	ug/l	1	-	U	Yes
2-Methylnaphthalene	0.96	ug/l	1	-	U	Yes
Naphthalene	0.96	ug/l	1	-	U	Yes

Sample ID: FA42192-15
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: AQ - Field Blank Water

METHOD: 8270D

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Caprolactam	10	ug/l	1	-	U	Yes
4-Chloro-3-methyl phenol	5.0	ug/l	1	-	U	Yes
2-Chlorophenol	5.0	ug/l	1	-	U	Yes
2,4-Dichlorophenol	5.0	ug/l	1	-	U	Yes
2,4-Dimethylphenol	5.0	ug/l	1	-	U	Yes
2,4-Dinitrophenol	25	ug/l	1	-	U	Yes
4,6-Dinitro-o-cresol	10	ug/l	1	-	U	Yes
2-Methylphenol	5.0	ug/l	1	-	U	Yes
3&4-Methylphenol	5.0	ug/l	1	-	U	Yes
2-Nitrophenol	5.0	ug/l	1	-	U	Yes
4-Nitrophenol	25	ug/l	1	-	U	Yes
Pentachlorophenol	25	ug/l	1	-	U	Yes
Phenol	5.0	ug/l	1	-	U	Yes
2,3,4,6-Tetrachlorophenol	5.0	ug/l	1	-	U	Yes
2,4,5-Trichlorophenol	5.0	ug/l	1	-	U	Yes
2,4,6-Trichlorophenol	5.0	ug/l	1	-	U	Yes
Acenaphthene	5.0	ug/l	1	-	U	Yes
Acenaphthylene	5.0	ug/l	1	-	U	Yes
Acetophenone	5.0	ug/l	1	-	U	Yes
Anthracene	5.0	ug/l	1	-	U	Yes
Atrazine	5.0	ug/l	1	-	U	Yes
Benzaldehyde	25	ug/l	1	-	U	Yes
Benzo(a)anthracene	5.0	ug/l	1	-	U	Yes
Benzo(a)pyrene	5.0	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	5.0	ug/l	1	-	U	Yes
Benzo(g,h,i)perylene	5.0	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	5.0	ug/l	1	-	U	Yes
1,1'-Biphenyl	5.0	ug/l	1	-	U	Yes
4-Bromophenyl phenyl ether	5.0	ug/l	1	-	U	Yes
Butyl benzyl phthalate	5.0	ug/l	1	-	U	Yes
Carbazole	5.0	ug/l	1	-	U	Yes
4-Chloroaniline	5.0	ug/l	1	-	U	Yes
bis(2-Chloroethoxy)methane	5.0	ug/l	1	-	UJ	Yes
bis(2-Chloroethyl)ether	5.0	ug/l	1	-	U	Yes
bis(2-Chloroisopropyl)ether	5.0	ug/l	1	-	U	Yes

2-Chloronaphthalene	5.0	ug/l	1	-	U	Yes
4-Chlorophenyl phenyl ether	5.0	ug/l	1	-	U	Yes
Chrysene	5.0	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	5.0	ug/l	1	-	U	Yes
Dibenzofuran	5.0	ug/l	1	-	U	Yes
3,3'-Dichlorobenzidine	5.0	ug/l	1	-	U	Yes
Diethyl phthalate	5.0	ug/l	1	-	U	Yes
Dimethyl phthalate	5.0	ug/l	1	-	U	Yes
Di-n-butyl phthalate	5.0	ug/l	1	-	U	Yes
Di-n-octyl phthalate	5.0	ug/l	1	-	U	Yes
2,4-Dinitrotoluene	5.0	ug/l	1	-	U	Yes
2,6-Dinitrotoluene	5.0	ug/l	1	-	U	Yes
bis(2-Ethylhexyl)phthalate	5.0	ug/l	1	-	U	Yes
Fluoranthene	5.0	ug/l	1	-	U	Yes
Fluorene	5.0	ug/l	1	-	U	Yes
Hexachlorobenzene	5.0	ug/l	1	-	U	Yes
Hexachlorobutadiene	5.0	ug/l	1	-	U	Yes
Hexachlorocyclopentadiene	5.0	ug/l	1	-	U	Yes
Hexachloroethane	5.0	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	5.0	ug/l	1	-	U	Yes
Isophorone	5.0	ug/l	1	-	U	Yes
2-Methylnaphthalene	5.0	ug/l	1	-	UJ	Yes
2-Nitroaniline	5.0	ug/l	1	-	U	Yes
3-Nitroaniline	5.0	ug/l	1	-	U	Yes
4-Nitroaniline	5.0	ug/l	1	-	U	Yes
Nitrobenzene	5.0	ug/l	1	-	U	Yes
N-Nitroso-di-n-propylamine	5.0	ug/l	1	-	U	Yes
N-Nitrosodiphenylamine	5.0	ug/l	1	-	U	Yes
Phenanthrene	5.0	ug/l	1	-	U	Yes
Pyrene	5.0	ug/l	1	-	U	Yes
1,2,4,5-Tetrachlorobenzene	5.0	ug/l	1	-	U	Yes
METHOD: 8270D (SIM)						
Benzo(a)anthracene	0.20	ug/l	1	-	U	Yes
Benzo(a)pyrene	0.20	ug/l	1	-	U	Yes
Benzo(b)fluoranthene	0.20	ug/l	1	-	U	Yes
Benzo(k)fluoranthene	0.20	ug/l	1	-	U	Yes
Chrysene	0.20	ug/l	1	-	U	Yes
Dibenzo(a,h)anthracene	0.20	ug/l	1	-	U	Yes
1,4-Dioxane	0.30	ug/l	1	-	U	Yes
Indeno(1,2,3-cd)pyrene	0.20	ug/l	1	-	U	Yes
1-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
2-Methylnaphthalene	1.0	ug/l	1	-	U	Yes
Naphthalene	1.0	ug/l	1	-	U	Yes

DATA REVIEW WORKSHEETS

Project Number:_FA42192_____
Date:_____March_16-17,_2017_____
Shipping Date:__March_17,_2017_____
EPA Region:_____2_____

REVIEW OF SEMIVOLATILE ORGANIC PACKAGE

The following guidelines for evaluating volatile organics were created to delineate required validation actions. This document will assist the reviewer in using professional judgment to make more informed decision and in better serving the needs of the data users. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence: EPA Hazardous Waste Support Section, SOP HW-35A, July 2015 –Revision 0. *Semivolatile Data Validation*. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

The hardcopied (laboratory name) _Accutest_____ data package received has been reviewed and the quality control and performance data summarized. The data review for SVOCs included:

Lab. Project/SDG No.: ____FA42192_____ Sample matrix: ___Groundwater____
No. of Samples: _____11SCAN/22SIM_____

Trip blank No.: _____-_____
Field blank No.: _____FA42192-1;_ FA42192-15_____
Equipment blank No.:_____-_____
Field duplicate No.:_____-_____

___X___ Data Completeness
___X___ Holding Times
___X___ GC/MS Tuning
___X___ Internal Standard Performance
___X___ Blanks
___X___ Surrogate Recoveries
___X___ Matrix Spike/Matrix Spike Duplicate
___X___ Laboratory Control Spikes
___X___ Field Duplicates
___X___ Calibrations
___X___ Compound Identifications
___X___ Compound Quantitation
___X___ Quantitation Limits

_Overall Comments:_Selected_SVOCs_from_the_TCL_special_list_analyzed_by_method_SW846-_8270D;_Selected_PAHs_and_1,4-Dioxane_analyzed_by_method_SW846-8270D_(SIM);_1,4-dioxane_and_PAH's_analyzed_separately._____

Definition of Qualifiers:

J- Estimated results
U- Compound not detected
R- Rejected data
UJ- Estimated nondetect

Reviewer:_Rafael DeLaut_____
Date:___April_24,_2017_____

DATA REVIEW WORKSHEETS

DATA COMPLETENESS

MISSING INFORMATION	DATE LAB. CONTACTED	DATE RECEIVED

All criteria were met _____
Criteria were not met
and/or see below ___X___

HOLDING TIMES

The objective of this parameter is to ascertain the validity of the results based on the holding time of the sample from time of collection to the time of analysis.

Complete table for all samples and note the analysis and/or preservation not within criteria

SAMPLE ID	DATE SAMPLED	DATE EXTRACTED/ANALYZED	pH	ACTION
FA42192-12	03/17/17	04/03/17	-	No action; confirmation run. Surrogate recovery outside laboratory control limits.
FA42192-13	03/17/17	04/03/17	-	
FA42192-14	03/17/17	04/03/17	-	
All samples extracted and analyzed within method recommended holding time except for the cases described in this document. Sample preservation appropriate.				

Cooler temperature (Criteria: 4 ± 2 °C): _____1.6_-_5.8_°C_____

Actions

Results will be qualified based on the criteria of the following Table:

Table 1. Holding Time Actions for Semivolatile Analyses

Matrix	Preserved	Criteria	Action: Detected Associated Compounds	Action: Non-Detected Associated Compounds
Aqueous	No	≤ 7 days (for extraction) ≤ 40 days (for analysis)	Use professional judgment	
	No	> 7 days (for extraction) > 40 days (for analysis)	J	Use professional judgment
	Yes	≤ 7 days (for extraction) ≤ 40 days (for analysis)	No qualification	
	Yes	> 7 days (for extraction) > 40 days (for analysis)	J	UJ
	Yes/No	Grossly Exceeded	J	UJ or R
Non-Aqueous	No	≤ 14 days (for extraction) ≤ 40 days (for analysis)	Use professional judgment	
	No	> 14 days (for extraction) > 40 days (for analysis)	J	Use professional judgment
	Yes	≤ 14 days (for extraction) ≤ 40 days (for analysis)	No qualification	
	Yes	> 14 days (for extraction) > 40 days (for analysis)	J	UJ
	Yes/No	Grossly Exceeded	J	UJ or R

All criteria were met _X__
Criteria were not met see below ____

GC/MS TUNING

The assessment of the tuning results is to determine if the sample instrumentation is within the standard tuning QC limits

_X__ The DFTPP performance results were reviewed and found to be within the specified criteria.

_X__ DFTPP tuning was performed for every 12 hours of sample analysis.

If no, use professional judgment to determine whether the associated data should be accepted, qualified or rejected.

Notes: These requirements do not apply when samples are analyzed by the Selected Ion Monitoring (SIM) technique.

All mass spectrometer conditions must be identical to those used during the sample analysis. Background subtraction actions resulting in spectral distortion are unacceptable

Notes: No data should be qualified based of DFTPP failure.

The requirement to analyze the instrument performance check solution is optional when analysis of PAHs/pentachlorophenol is to be performed by the SIM technique.

List the samples affected:
_____
_____
_____
_____

Actions:

1. If sample are analyzed without a preceding valid instrument performance check or are analyzed 12 hours after the Instrument Performance Check, qualify all data in those samples as unusable (R).
2. If ion abundance criteria are not met, use professional judgment to determine to what extent the data may be utilized.
3. State in the Data Review Narrative, decisions to use analytical data associated with DFTPP instrument performance checks not meeting the contract requirements.
4. Use professional judgment to determine if associated data should be qualified based on the spectrum of the mass calibration compounds.

All criteria were met ___X___
Criteria were not met
and/or see below _____

INITIAL CALIBRATION VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:____03/26/17_(SCAN)_____ _____02/13/17;_04/03/17_(SIM)_____
Instrument ID numbers:_____GCMSL_____ _____GCMSW_____
Matrix/Level:_____Aqueous/low_____ _____Aqueous/low_____

Date of initial calibration:___02/23/17_(SIM)_____
Instrument ID numbers:_____GCMSU_____
Matrix/Level:_____Aqueous/low_____

DATE	LAB FILE ID#	CRITERIA OUT RFs, %RSD, %D, r	COMPOUND	SAMPLES AFFECTED
Initial calibration meets the method and guidance validation document performance criteria.				

Note:

Actions:

Qualify the initial calibration analytes listed in Table 2 using the following criteria:

Table 3. Initial Calibration Actions for Semivolatile Analysis

Criteria	Action	
	Detect	Non-detect
Initial Calibration not performed at specified frequency and sequence	Use professional judgment R	Use professional judgment R
Initial Calibration not performed at the specified concentrations	J	UJ
RRF < Minimum RRF in Table 2 for target analyte	Use professional judgment J+ or R	R
RRF ≥ Minimum RRF in Table 2 for target analyte	No qualification	No qualification
%RSD > Maximum %RSD in Table 2 for target analyte	J	Use professional judgment
%RSD ≤ Maximum %RSD in Table 2 for target analyte	No qualification	No qualification

Initial Calibration

Table 2. RRF, %RSD, and %D Acceptance Criteria in Initial Calibration and CCV for Semivolatil Analysis

Analyte	Minimum RRF	Maximum %RSD	Opening Maximum %D[1]	Opening Maximum %D[1]
1,4-Dioxane	0.010	40.0	± 40.0	± 50.0
Benzaldehyde	0.100	40.0	± 40.0	± 50.0
Phenol	0.080	20.0	± 20.0	± 25.0
Bis(2-chloroethyl)ether	0.100	20.0	± 20.0	± 25.0
2-Chlorophenol	0.200	20.0	± 20.0	± 25.0
2-Methylphenol	0.010	20.0	± 20.0	± 25.0
3-Methylphenol	0.010	20.0	± 20.0	± 25.0
2,2'-Oxybis-(1-chloropropane)	0.010	20.0	± 25.0	± 50.0
Acetophenone	0.060	20.0	± 20.0	± 25.0
4-Methylphenol	0.010	20.0	± 20.0	± 25.0
N-Nitroso-di-n-propylamine	0.080	20.0	± 25.0	± 25.0
Hexachloroethane	0.100	20.0	± 20.0	± 25.0
Nitrobenzene	0.090	20.0	± 20.0	± 25.0
Isophorone	0.100	20.0	± 20.0	± 25.0
2-Nitrophenol	0.060	20.0	± 20.0	± 25.0
2,4-Dimethylphenol	0.050	20.0	± 25.0	± 50.0
Bis(2-chloroethoxy)methane	0.080	20.0	± 20.0	± 25.0
2,4-Dichlorophenol	0.060	20.0	± 20.0	± 25.0
Naphthalene	0.200	20.0	± 20.0	± 25.0
4-Chloroaniline	0.010	40.0	± 40.0	± 50.0
Hexachlorobutadiene	0.040	20.0	± 20.0	± 25.0
Caprolactam	0.010	40.0	± 30.0	± 50.0
4-Chloro-3-methylphenol	0.040	20.0	± 20.0	± 25.0
2-Methylnaphthalene	0.100	20.0	± 20.0	± 25.0
Hexachlorocyclopentadiene	0.010	40.0	± 40.0	± 50.0
2,4,6-Trichlorophenol	0.090	20.0	± 20.0	± 25.0
2,4,5-Trichlorophenol	0.100	20.0	± 20.0	± 25.0
1,1'-Biphenyl	0.200	20.0	± 20.0	± 25.0

+

Analyte	Minimum RRF	Maximum %RSD	Opening Maximum %D[1]	Opening Maximum %D[1]
2-Chloronaphthalene	0.300	20.0	± 20.0	± 25.0
2-Nitroaniline	0.060	20.0	± 25.0	± 25.0
Dimethylphthalate	0.300	20.0	± 25.0	± 25.0
2,6-Dinitrotoluene	0.080	20.0	± 20.0	± 25.0
Acenaphthylene	0.400	20.0	± 20.0	± 25.0
3-Nitroaniline	0.010	20.0	± 25.0	± 50.0
Acenaphthene	0.200	20.0	± 20.0	± 25.0
2,4-Dinitrophenol	0.010	40.0	± 50.0	± 50.0
4-Nitrophenol	0.010	40.0	± 40.0	± 50.0
Dibenzofuran	0.300	20.0	± 20.0	± 25.0
2,4-Dinitrotoluene	0.070	20.0	± 20.0	± 25.0
Diethylphthalate	0.300	20.0	± 20.0	± 25.0
1,2,4,5-Tetrachlorobenzene	0.100	20.0	± 20.0	± 25.0
4-Chlorophenyl-phenylether	0.100	20.0	± 20.0	± 25.0
Fluorene	0.200	20.0	± 20.0	± 25.0
4-Nitroaniline	0.010	40.0	± 40.0	± 50.0
4,6-Dinitro-2-methylphenol	0.010	40.0	± 30.0	± 50.0
4-Bromophenyl-phenyl ether	0.070	20.0	± 20.0	± 25.0
N-Nitrosodiphenylamine	0.100	20.0	± 20.0	± 25.0
Hexachlorobenzene	0.050	20.0	± 20.0	± 25.0
Atrazine	0.010	40.0	± 25.0	± 50.0
Pentachlorophenol	0.010	40.0	± 40.0	± 50.0
Phenanthrene	0.200	20.0	± 20.0	± 25.0
Anthracene	0.200	20.0	± 20.0	± 25.0
Carbazole	0.050	20.0	± 20.0	± 25.0
Di-n-butylphthalate	0.500	20.0	± 20.0	± 25.0
Fluoranthene	0.100	20.0	± 20.0	± 25.0
Pyrene	0.400	20.0	± 25.0	± 50.0
Butylbenzylphthalate	0.100	20.0	± 25.0	± 50.0

Analyte	Minimum RRF	Maximum %RSD	Opening Maximum %D[1]	Opening Maximum %D[1]
3,3'-Dichlorobenzidine	0.010	40.0	± 40.0	± 50.0
Benzo(a)anthracene	0.300	20.0	± 20.0	± 25.0
Chrysene	0.200	20.0	± 20.0	± 50.0
Bis(2-ethylhexyl) phthalate	0.200	20.0	± 25.0	± 50.0
Di-n-octylphthalate	0.010	40.0	± 40.0	± 50.0
Benzo(b)fluoranthene	0.010	20.0	± 25.0	± 50.0
Benzo(k)fluoranthene	0.010	20.0	± 25.0	± 50.0
Benzo(a)pyrene	0.010	20.0	± 20.0	± 50.0
Indeno(1,2,3-cd)pyrene	0.010	20.0	± 25.0	± 50.0
Dibenzo(a,h)anthracene	0.010	20.0	± 25.0	± 50.0
Benzo(g,h,i)perylene	0.010	20.0	± 30.0	± 50.0
2,3,4,6-Tetrachlorophenol	0.040	20.0	± 20.0	± 50.0
Naphthalene	0.600	20.0	± 25.0	± 25.0
2-Methylnaphthalene	0.300	20.0	± 20.0	± 25.0
Acenaphthylene	0.900	20.0	± 20.0	± 25.0
Acenaphthene	0.500	20.0	± 20.0	± 25.0
Fluorene	0.700	20.0	± 25.0	± 50.0
Phenanthrene	0.300	20.0	± 25.0	± 50.0
Anthracene	0.400	20.0	± 25.0	± 50.0
Fluoranthene	0.400	20.0	± 25.0	± 50.0
Pyrene	0.500	20.0	± 30.0	± 50.0
Benzo(a)anthracene	0.400	20.0	± 25.0	± 50.0
Chyrsene	0.400	20.0	± 25.0	± 50.0
Benzo(b)fluoranthene	0.100	20.0	± 30.0	± 50.0
Benzo(k)fluoranthene	0.100	20.0	± 30.0	± 50.0
Benzo(a)pyrene	0.100	20.0	± 25.0	± 50.0
Indeno(1,2,3-cd)pyrene	0.100	20.0	± 40.0	± 50.0
Dibenzo(a,h)anthracene	0.010	25.0	± 40.0	± 50.0
Benzo(g,h,i)perylene	0.020	25.0	± 40.0	± 50.0

Pentachlorophenol	0.010	40.0	± 50.0	± 50.0
Deuterated Monitoring Compounds				
Analyte	**Minimum RRF**	**Maximum %RSD**	**Opening Maximum %D[1]**	**Closing Maximum %D**
1,4-Dioxane-d_8	0.010	20.0	± 25.0	± 50.0
Phenol-d_5	0.010	20.0	± 25.0	± 25.0
Bis-(2-chloroethyl)ether-d_8	0.100	20.0	± 20.0	± 25.0
2-Chlorophenol-d_4	0.200	20.0	± 20.0	± 25.0
4-Methylphenol-d_8	0.010	20.0	± 20.0	± 25.0
4-Chloroaniline-d_4	0.010	40.0	± 40.0	± 50.0
Nitrobenzene-d_5	0.050	20.0	± 20.0	± 25.0
2-Nitrophenol-d_4	0.050	20.0	± 20.0	± 25.0
2,4-Dichlorophenol-d_3	0.060	20.0	± 20.0	± 25.0
Dimethylphthalate-d_6	0.300	20.0	± 20.0	± 25.0
Acenaphthylene-d_8	0.400	20.0	± 20.0	± 25.0
4-Nitrophenol-d_4	0.010	40.0	± 40.0	± 50.0
Fluorene-d_{10}	0.100	20.0	± 20.0	± 25.0
4,6-Dinitro-2-methylphenol-d_2	0.010	40.0	± 30.0	± 50.0
Anthracene-d_{10}	0.300	20.0	± 20.0	± 25.0
Pyrene-d_{10}	0.300	20.0	± 25.0	± 50.0
Benzo(a)pyrene-d_{12}	0.010	20.0	± 20.0	± 50.0
Fluoranthene-d_{10} (SIM)	0.400	20.0	± 25.0	± 50.0
2-Methylnaphthalene-d_{10} (SIM)	0.300	20.0	± 20.0	± 25.0

[1] If a closing CCV is acting as an opening CCV, all target analytes must meet the requirements for an opening CCV.

Note: If analysis by SIM technique is requested for PAH/pentachlorophenols, calibration standards analyzed at 0.10, 0.20, 0.40, 0.80, and 1.0 ng/uL for each target compound of interest and the associated DMCs. Pentachlorophenol will require only a four point initial calibration at 0.20, 0.40, 0.80, and 1.0 ng/uL.

All criteria were met __X_____
Criteria were not met
and/or see below _____

CONTINUING CALIBRATION VERIFICATION/

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:_____02/13/17;_04/03/17_(SIM)_____ ____02/23/17_(SIM)_____
Date of initial calibration verification (ICV):_02/13/17;_04/03/17_____ ____02/23/17_____
Date of continuing calibration verification (CCV):_03/24/17;_03/28/17___ _03/24/17;_03/25/17_____
_03/31/17_____ _03/29/17_____
Date of closing CCV:__02/14/17;_03/21/17;_04/01/17;_04/03/17_____ _____03/29/17_____
Instrument ID numbers:_____GCMSW_____ _____GCMSU_____
Matrix/Level:_____Aqueous/low_____ _____Aqueous/low_____

Date of initial calibration:_____03/26/17_(SCAN)_____
Date of initial calibration verification (ICV):__03/2617_____
Date of continuing verification (CCV):_____03/27/17;_03/28/17;_03/30/17;_03/31/17_____
Date of closing CCV:_____-_____
Instrument ID numbers:_____GCMSL_____
Matrix/Level:_____Aqueous/low_____

DATE	LAB FILE ID#	CRITERIA OUT RFs, %RSD, %D, r	COMPOUND	SAMPLES AFFECTED
GCMSL				
03/26/17	icv3976-50, 2x	29.3	bis(2-chloroethoxy) methane	FA42192-8; -11; -12; -13; -14; -15
	icv3976-50, benz	22.5	2-methylnaphthalene	
03/30/17	cc3976-50	-29.9	2-nitrophenol	FA42192-1; -2; -3; -7
		-21.0	Hexachlorocyclopentadiene^	
		-27.8	2-nitroaniline	
		-24.7	3-nitroaniline^	
		-45.6	2,4-dinitrophenol^	
		-24.6	2,4-dinitrotoluene	
		-21.0	2,3,4,6-tetrachlorophenol	
		-47.9	4,6-dinitro-o-cresol	
03/31/17	cc3976-50	-44.5	2-nitrophenol	FA42192-6
		-34.6	Hexachlorocyclopentadiene^	
		-31.4	2-nitroaniline	
		-90.7	2,4-dinitrophenol^	
		-33.6	2,4-dinitrotoluene	
		-21.3	2,3,4,6-tetrachlorophenol	
		-78.4	4,6-dinitro-o-cresol	
		-35.0	Pentachlorophenol^	

Note: Initial and continuing calibration verifications meet the method and guidance document required performance criteria except for the cases described in this document.

^Analytes with % difference outside the method performance criteria but within guidance document performance criteria in the continuing calibration verification guidance document.

No closing calibration verification included in data package for instrument GCMSL (SCAN). No action taken, professional judgment.

Analytes not meeting the initial and continuing % difference in the initial and continuing calibration verification criteria qualified as estimated (J or UJ) in affected samples.

Actions:

Notes: Verify that the CCV is run at the required frequency (an opening and closing CCV must be run within 12-hour period).

All DMCs must meet the RRF values given in Table 2. No qualification of the data is necessary on DMCs RRF and %RSD/%D alone. Use professional judgment to evaluate DMCs and %RSD/%D data in conjunction with DMCs recoveries to determine the need for qualification of the data.

Qualify the initial calibration analytes listed in Table 2 using the following criteria in the CCVs:

Table 4. CCV Actions for Semivolatile Analysis

Criteria for Opening CCV	Criteria for Closing CCV	Action	
		Detect	Non-detect
CCV not performed at required frequency and sequence	CCV not performed at required frequency	Use professional judgment R	Use professional judgment R
CCV not performed at specified concentration	CCV not performed at specified concentration	Use professional judgment	Use professional judgment
RRF < Minimum RRF in Table 2 for target analyte	RRF < Minimum RRF in Table 2 for target analyte	Use professional judgment J or R	R
RRF ≥ Minimum RRF in Table 2 for target analyte	RRF ≥ Minimum RRF in Table 2 for target analyte	No qualification	No qualification
%D outside the Opening Maximum %D limits in Table 2 for target analyte	%D outside the Closing Maximum %D limits in Table 2 for target analyte	J	UJ
%D within the inclusive Opening Maximum %D limits in Table 2 for target analyte	%D within the inclusive Closing Maximum %D limits in Table 2 for target analyte	No qualification	No qualification

All criteria were met _____
Criteria were not met
and/or see below ____X____

BLANK ANALYSIS RESULTS (Sections 1 & 2)

The assessment of the blank analysis results is to determine the existence and magnitude of contamination problems. The criteria for evaluation of blanks apply only to blanks associated with the samples, including trip, equipment, and laboratory blanks. If problems with any blanks exist, all data associated with the case must be carefully evaluated to determine whether or not there is an inherent variability in the data for the case, or if the problem is an isolated occurrence not affecting other data.

List the contamination in the blanks below. High and low levels blanks must be treated separately.

Notes: The concentration of non-target compounds in all blanks must be less than or equal to 10 ug/L.

The concentration of target compounds in all blanks must be less than its CRQL listed in the method.

Samples taken from a drinking water tap do not have and associated field blank.

Laboratory blanks

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
_No_target_analytes_detected_in_method_blanks_except_for_the_cases_described_in_this_document				
_03/25/17_____	_OP64320-MB__	__Aq./low___	_1,4-Dioxane_____	__0.49_ug/l_____

Note: Laboratory qualified the results in samples FA42192-8, FA42192-11, FA42192-12, FA42192-13, FA42192-14, and FA42192-15 with a B qualifier.

Validator qualification based on guidelines: - Positive results below the reporting limits are reported at the reporting limits and qualified as non-detected (U). Results above the reporting limits and below method blanks results are reported at sample results and qualified as non-detects (U). Results above the reporting limits and blank result are not qualified, professional judgment.

Field/Equipment/Trip blank

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
_No_target_analytes_detected_in_the_field_blanks_analyzed_with_this_data_package._No_equipment _blank_analyzed_with_this_data_package._____				

Note:

All criteria were met __X__
Criteria were not met
and/or see below _____

BLANK ANALYSIS RESULTS (Section 3)

Blank Actions

Qualify samples based on the criteria summarized in Table 5:

Table 5. Blank and TCLP/SPLP LEB Actions for Semivolatile Analysis

Blank Type	Blank Result	Sample Result	Action
Method, TCLP/SPLP LEB, Field	Detect	Non-detect	No qualification
	< CRQL	< CRQL	Report at CRQL and qualify as non-detect (U)
		≥ CRQL	Use professional judgment
	≥ CRQL	< CRQL	Report at CRQL and qualify as non-detect (U)
		≥ CRQL but < Blank Result	Report at sample results and qualify as non-detect (U) or as unusable (R)
		≥ CRQL and ≥ Blank Result	Use professional judgment
	Grossly high	Detect	Report at sample results and qualify as unusable (R)
	TIC > 5.0 ug/L (water) or 0.0050 mg/L (TCLP leachate) or TIC > 170 ug/Kg (soil)	Detect	Use professional judgment

List samples qualified

CONTAMINATION SOURCE/LEVEL	COMPOUND	CONC/UNITS	AL/UNITS	SQL	AFFECTED SAMPLES

All criteria were met ____
Criteria were not met
and/or see below ___X___

SURROGATE SPIKE RECOVERIES – DEUTERATED MONITORING COMPOUNDS (DMCs)

Laboratory performance of individual samples is established by evaluation of surrogate spike recoveries – deuterated monitoring compounds. All samples are spiked with surrogate compounds prior to sample analysis. The accuracy of the analysis is measured by the surrogate percent recovery. Since the effects of the sample matrix are frequently outside the control of the laboratory and may present relatively unique problems, the validation of data is frequently subjective and demands analytical experience and professional judgment.

Notes: Recoveries for DMCs in samples and blanks must be within the limits specified in Table 7.

The recovery limits for any of the compounds listed in Table 7 may be expanded at any time during the period of performance if USEPA determines that the limits are too restrictive.

If a DMC is not added in the samples and blanks or the concentrations of DMCs in the samples and blank not the specified, use professional judgment in qualifying the data.

Table 7. DMC Actions for Semivolatile Analysis

Criteria	Action	
	Detect	Non-detect
%R < 10% (excluding DMCs with 10% as a lower acceptance limit)	J-	R
10% ≤ %R (excluding DMCs with 10% as a lower acceptance limit) < Lower Acceptance Limit	J-	UJ
Lower Acceptance limit ≤ %R ≤ Upper Acceptance Limit	No qualification	No qualification
%R > Upper Acceptance Limit	J+	No qualification

List the percent recoveries (%Rs) which do not meet the criteria for DMCs (surrogate) recovery.

Matrix:___Groundwater_____

SAMPLE ID **SURROGATE COMPOUND** **ACTION**

_DMCs_meet_the_required_criteria_in_all_samples_analyzed_except_for_the_cases_described_in___
_this_document._Non-_deuterated_surrogates_added_to_the_samples_and_were_within_laboratory_
_recovery_limits_except_in_the_cases_described_in_this_document._____
_FA42192-6_____2-Fluorophenol_5_%_(14_-_67_control_limit)_____
_____2,4,6-tribromophenol_122_%_(33_-_118_control_limit)_____

Note: Surrogate standards not recovered in sample FA42192-13 (SIM) due to dilution. No action taken, professional judgment. Surrogates recovered in the undiluted sample.

No qualification made on sample FA42192-6, professional judgment.

Terphenyl-d14 recovered low in samples FA42192-12 (31 %); FA42192-13 (32 %); and FA42192-14 (29 %); control limits 39 – 121 %. Confirmed by re-extraction and analysis. Samples extracted beyond holding time. No action taken, professional judgment.

Surrogate recoveries in samples analyzed by SW846-8270D (SIM) were corrected for the concentration of the surrogate spiked. The correction is due to the following: the laboratory split batches between 8270 and 8270SIM. The surrogates were spiked at 8270 levels, which is 5x higher than what the laboratory 8270SIM setup expects. SIM surrogates recoveries were divided by 5. Calculations were spot checked.

Table 8. Semivolatile DMCs and the Associated Target Analytes

1,4–Dioxane-d_8 (DMC-1)	**Phenol-d_5 (DMC-2)**	**Bis(2-Chloroethyl) ether-d_8 (DMC-3)**
1,4-Dioxane	Benzaldehyde Phenol	Bis(2-chloroethyl)ether 2,2'-Oxybis(1-chloropropane) Bis(2-chloroethoxy)methane
2-Chlorophenol-d_4 (DMC-4)	**4-Methylphenol-d_8 (DMC-5)**	**4-Chloroaniline-d_4 (DMC-6)**
2-Chlorophenol	2-Methylphenol 3-Methylphenol 4-Methylphenol 2,4-Dimethylphenol	4-Chloroaniline Hexachlorocyclopentadiene Dichlorobenzidine
Nitrobenzene-d_5 (DMC-7)	**2-Nitrophenol-d_4 (DMC-8)**	**2,4-Dichlorophenol-d_3 (DMC-9)**
Acetophenone N-Nitroso-di-n-propylamine Hexachloroethane Nitrobenzene 2,6-Dinitrotoluene 2,4-Dinitrotoluene N-Nitrosodiphenylamine	Isophorone 2-Nitrophenol	2,4-Dichlorophenol Hexachlorobutadiene Hexachlorocyclopentadiene 4-Chloro-3-methylphenol 2,4,6-Trichlorophenol 2,4,5-Trichlorophenol 1,2,4,5-Tetrachlorobenzene *Pentachlorophenol 2,3,4,6-Tetrachlorophenol
Dimethylphthalate-d_6 (DMC-10)	**Acenaphthylene-d_8 (DMC-11)**	**4-Nitrophenol-d_4 (DMC-12)**
Caprolactam 1,1'-Biphenyl Dimethylphthalate Diethylphthalate Di-n-butylphthalate Butylbenzylphthalate Bis(2-ethylhexyl) phthalate Di-n-octylphthalate	*Naphthalene *2-Methylnaphthalene 2-Chloronaphthalene *Acenaphthylene *Acenaphthene	2-Nitroaniline 3-Nitroaniline 2,4-Dinitrophenol 4-Nitrophenol 4-Nitroaniline

Fluorene-d_{10} (DMC-13)	**4,6-Dinitro-2-methylphenol-d_2 (DMC-14)**	**Anthracene-d_{10} (DMC-15)**
Dibenzofuran *Fluorene 4-Chlorophenyl-phenylether 4-Bromophenyl-phenylether Carbazole	4,6-Dinitro-2-methylphenol	Hexachlorobenzene Atrazine *Phenanthrene *Anthracene
Pyrene-d_{10} (DMC-16)	**Benzo(a)pyrene-d_{12} (DMC-17)**	
*Fluoranthene *Pyrene *Benzo(a)anthracene *Chrysene	3,3'-Dichlorobenzidine *Benzo(b)fluoranthene *Benzo(k)fluoranthene *Benzo(a)pyrene *Indeno(1,2,3-cd)pyrene *Dibenzo(a,h)anthracene *Benzo(g,h,i)perylene	

*Included in optional Target Analyte List (TAL) of PAHs and PCP only.

Table 9. Semivolatile SIM DMCs and the Associated Target Analytes

Fluoranthene-d10 (DMC-1)	**2-Methylnaphthalene-d10 (DMC-2)**
Fluoranthene	Naphthalene
Pyrene	2-Methylnaphthalene
Benzo(a)anthracene	Acenaphthylene
Chrysene	Acenaphthene
Benzo(b)fluoranthene	Fluorene
Benzo(k)fluoranthene	Pentachlorophenol
Benzo(a)pyrene	Phenanthrene
Indeno(1,2,3-cd)pyrene	Anthracene
Dibenzo(a,h)anthracene	
Benzo(g,h,i)perylene	

All criteria were met _____
Criteria were not met
and/or see below ____X___

VII. A MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

This data is generated to determine long term precision and accuracy in the analytical method for various matrices. This data alone cannot be used to evaluate the precision and accuracy of individual samples. If any % R in the MS or MSD falls outside the designated range, the reviewer should determine if there are matrix effects, i.e. LCS data are within the QC limits but MS/MSD data are outside QC limit.

1. MS/MSD Recoveries and Precision Criteria

The laboratory should use one MS and a duplicate analysis of an unspiked field sample if target analytes are expected in the sample. If target analytes are not expected, MS/MSD should be analyzed.

NOTES: Data for MS and MSDs will not be present unless requested by the Region. Notify the Contract Laboratory COR if a field or trip blank was used for the MS and MSD.

For a Matrix Spike that does not meet criteria, apply the action to only the field sample used to prepare the Matrix Spike sample. If it is clearly stated in the data validation materials that the samples were taken through incremental sampling or some other method guaranteeing the homogeneity of the sample group, then the entire sample group may be qualified.

List the %Rs, RPD of the compounds which do not meet the criteria.

Sample ID:_____FA42192-6_____ Matrix/Level:__Groundwater_
Sample ID:_____FA42192-13_____ Matrix/Level:__Groundwater_
Sample ID:_____JC42157-7_(SIM)_____ Matrix/Level:__Groundwater_
Sample ID:_____JC42192-15_(SIM)_____ Matrix/Level:__Groundwater_

The QC reported here applies to the following samples: **Method:** SW846 8270D
FA42192-1, FA42192-2, FA42192-3, FA42192-6, FA42192-7

Compound	FA42192-6 ug/l	Q	Spike ug/l	MS ug/l	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
2,4-Dinitrophenol	ND		192	212	110	192	239	124*	12	44-112/25
4,6-Dinitro-o-cresol	ND		192	270	140*	192	292	152*	8	66-121/23
2-Nitrophenol	ND		96.2	114	119*	96.2	105	109*	8	53-102/29
Pentachloro-phenol	ND		192	259	135*	192	237	123*	9	61-115/26
Phenol	ND		96.2	55.2	57*	96.2	44.4	46	22	19-56/35
2,4-Dinitrotoluene	ND		96.2	110	114*	96.2	107	111*	3	61-110/21
Isophorone	ND		96.2	87.8	91*	96.2	78.1	81	12	43-87/25
3-Nitroaniline	ND		96.2	106	110*	96.2	95.4	99	11	56-106/27
Nitrobenzene	ND		96.2	102	106*	96.2	91.6	95	11	50-104/28

* - outside laboratory control limits

Note: As per chain-of-custody form sample FA42192-3 was to use as the QC sample (MS/MSD) for this batch. Samples FA42192-6; FA42192-13; FA42157-7; and FA42237-15 were used instead. No action taken, professional judgment.

MS/MSD % recovery and RPD within laboratory control limits except for the cases described in this document.

Results for samples having either MS or the MSD outside the laboratory control limits are not qualified, professional judgment.

Samples with MS/MSD % recovery high are not qualified, professional judgment. 1,4-dioxane not qualified. No positive results associated with sample batch.

The QC reported here applies to the following samples: **Method:** SW846 8270D
FA42192-8, FA42192-11, FA42192-12, FA42192-13, FA42192-14, FA42192-15

Compound	FA42192-13 ug/l	Q	Spike ug/l	MS ug/l	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
bis(2-Chloroethoxy) methane	ND		96.2	48.3	50*	96.2	50.8	53	5	51-102/28

Note: MS/MSD % recovery and RPD within laboratory control limits except for the cases described in this document.

Results for samples having either MS or the MSD outside the laboratory control limits are not qualified, professional judgment.

Note:

* QC limits are laboratory in-house performance criteria, LL = lower limit, UL = upper limit.
* If QC limits are not available, use limits of 70 – 130 %.

Actions:

QUALITY	%R < LL	%R > UL
Positive results	J	J
Nondetects results	R	Accept

MS/MSD criteria apply only to the unspiked sample, its dilutions, and the associated MS/MSD samples:

If the % R for the affected compounds were < LL (or 70 %), qualify positive results (J) and nondetects (UJ).
If the % R for the affected compounds were > UL (or 130 %), only qualify positive results (J).
If 25 % or more of all MS/MSD %R were < LL (or 70 %) or if two or more MS/MSD %Rs were < 10%, qualify all positive results (J) and reject nondetects (R).

A separate worksheet should be used for each MS/MSD pair.

All criteria were met __X__
Criteria were not met
and/or see below _____

INTERNAL STANDARD PERFORMANCE

The assessment of the internal standard (IS) parameter is used to assist the data reviewer in determining the condition of the analytical instrumentation.

List the internal standard area of samples which do not meet the criteria.

DATE	SAMPLE ID	IS OUT	IS AREA	ACCEPTABLE RANGE	ACTION
Internal area meets the required criteria for batch samples corresponding to this data package.					

Action:

1. If an internal standard area count for a sample or blank is greater than 213.0% of the area for the associated standard (opening CCV or mid-point standard from initial calibration) (see Table 10 below):
   a. Qualify detects for compounds quantitated using that internal standard as estimated low (J-).
   b. Do not qualify non-detected associated compounds.
2. If an internal standard area count for a sample or blank is less than 20.0% of the area for the associated standard (opening CCV or mid-point standard from initial calibration):
   a. Qualify detects for compounds quantitated using that internal standard as estimated high (J+).
   b. Qualify non-detected associated compounds as unusable (R).
3. If an internal standard area count for a sample or blank is greater than or equal to 50.0%, and less than or equal to 213% of the area for the associated standard opening CCV or mid-point standard from initial calibration, no qualification of the data is necessary.
4. If an internal standard RT varies by more than 10.0 seconds: Examine the chromatographic profile for that sample to determine if any false positives or negatives exist. For shifts of a large magnitude, the reviewer may consider partial or total rejection of the data for that sample fraction. Detects should not need to be qualified as unusable (R) if the mass spectral criteria are met.
5. If an internal standard RT varies by less than or equal to 10.0 seconds, no qualification of the data is necessary.

Note: Inform the Contract Laboratory Program Project Officer (CLP PO) if the internal standard performance criteria are grossly exceeded. Note in the Data Review Narrative potential effects on the data resulting from unacceptable internal standard performance.

State in the Data Review Narrative if the required internal standard compounds are not added to a sample or blank or if the required internal standard compound is not analyzed at the specified concentration.

Actions:

Table 10. Internal Standard Actions for Semivolatile Analysis

Criteria	Action	
	Detect	Non-detect
Area response < 20% of the opening CCV or mid-point standard CS3 from ICAL	J+	R
20% < Area response < 50% of the opening CCV or mid-point standard CS3 from ICAL	J+	UJ
50% ≤ Area response ≤ 200% of the opening CCV or mid-point standard CS3 from ICAL	No qualification	No qualification
Area response > 200% of the opening CCV or mid-point standard CS3 from ICAL	J-	No qualification
RT shift between sample/blank and opening CCV or mid-point standard CS3 from ICAL > 10.0 seconds	R	R
RT shift between sample/blank and opening CCV or mid-point standard CS3 from ICAL < 10.0 seconds	No qualification	No qualification

All criteria were met __X__
Criteria were not met
and/or see below _____

TARGET COMPOUND IDENTIFICATION

Criteria:

Is the Relative Retention Times (RRTs) of reported compounds within ±0.06 RRT units of the standard RRT [opening Continuing Calibration Verification (CCV) or mid-point standard from the initial calibration]. **Yes?** or No?

List compounds not meeting the criteria described above:

Sample ID	Compounds	Actions
_____	_____	_____
_____	_____	_____
_____	_____	_____

Mass spectra of the sample compound and a current laboratory-generated standard [i.e., the mass spectrum from the associated calibration standard (opening CCV or mid-point standard from initial calibration)] must match according to the following criteria:

a. All ions present in the standard mass spectrum at a relative intensity greater than 10% must be present in the sample spectrum.
b. The relative intensities of these ions must agree within ±20% between the standard and sample spectra (e.g., for an ion with an abundance of 50% in the standard spectrum, the corresponding sample ion abundance must be between 30-70%).
c. Ions present at greater than 10% in the sample mass spectrum, but not present in the standard spectrum, must be evaluated by a reviewer experienced in mass spectral interpretation.

List compounds not meeting the criteria described above:

Sample ID	Compounds	Actions
_____	_____	_____
_Identified_compounds_meet_the_required_criteria____		_____
_____	_____	_____

Action:

1. The application of qualitative criteria for GC/MS analysis of target compounds requires professional judgment. It is up to the reviewer's discretion to obtain additional information from the laboratory. If it is determined that incorrect identifications were made, qualify all such data as unusable (R).

2. Use professional judgment to qualify the data if it is determined that cross-contamination has occurred.

3. Note in the Data Review Narrative any changes made to the reported compounds or concerns regarding target compound identifications. Note, for Contract Laboratory COR action, the necessity for numerous or significant changes.

TENTATIVELY IDENTIFIED COMPOUNDS (TICS)

NOTE: Tentatively identified compounds should only be evaluated when requested by a party from outside of the Hazardous Waste Support Section (HWSS).

List TICs

Sample ID	Compound	Sample ID	Compound

Action:

1. Qualify all TIC results for which there is presumptive evidence of a match (e.g. greater than or equal to 85% match) as tentatively identified (NJ), with approximated concentrations. TICs labeled "unknown" are qualified as estimated (J).
2. General actions related to the review of TIC results are as follows:
   a. If it is determined that a tentative identification of a non-target compound is unacceptable, change the tentative identification to "unknown" or another appropriate identification, and qualify the result as estimated (J).
   b. If all contractually-required peaks were not library searched and quantitated, the Region's designated representative may request these data from the laboratory.
3. In deciding whether a library search result for a TIC represents a reasonable identification, use professional judgment. If there is more than one possible match, report the result as "either compound X or compound Y". If there is a lack of isomer specificity, change the TIC result to a nonspecific isomer result (e.g., 1,3,5-trimethyl benzene to trimethyl benzene isomer) or to a compound class (e.g., 2-methyl, 3-ethyl benzene to a substituted aromatic compound).
4. The reviewer may elect to report all similar compounds as a total (e.g., all alkanes may be summarized and reported as total hydrocarbons).

5. Target compounds from other fractions and suspected laboratory contaminants should be marked as "non-reportable".
6. Other Case factors may influence TIC judgments. If a sample TIC match is poor, but other samples have a TIC with a valid library match, similar RRT, and the same ions, infer identification information from the other sample TIC results.
7. Note in the Data Review Narrative any changes made to the reported data or any concerns regarding TIC identifications.
8. Note, for Contract Laboratory COR action, failure to properly evaluate and report TICs

All criteria were met __X__
Criteria were not met
and/or see below _____

SAMPLE QUANTITATION AND REPORTED CONTRACT REQUIRED QUANTITATION LIMITS (CRQLS)

Action:

1. When a sample is analyzed at more than one dilution, the lower CRQL are used unless a QC exceedance dictates the use of higher CRQLs from the diluted sample. Samples reported with an "E" qualifier should be reported from the diluted sample.
2. If any discrepancies are found, the Region's designated representative may contact the laboratory to obtain additional information that could resolve any differences. If a discrepancy remains unresolved, the reviewer must use professional judgment to decide which value is the most accurate. Under these circumstances, the reviewer may determine that qualification of data is warranted. Note in the Data Review Narrative a description of the reasons for data qualification and the qualification that is applied to the data.
3. For non-aqueous samples, if the solids is less than 10.0%, use professional judgment for both detects and non-detects. If the percent solid for a soil sample is greater than or equal to 10.0% and less than 30.0%, use professional judgment to qualify detects and non-detects. If the percent solid for a soil sample is greater than or equal to 30.0%, detects and non-detects should not be qualified (see Table 11).
4. Note, for Contract Laboratory COR action, numerous or significant failures to accurately quantify the target compounds or to properly evaluate and adjust CRQLs.
5. Results between MDL and CRQL should be qualified as estimated "J".
6. Results < MDL should be reported at the CRQL and qualified "U". MDLs themselves should not be reported.

Table 11. Percent Solids Actions for Semivolatile Analysis for Non-Aqueous Samples

Criteria	Action	
	Detects	Non-detects
%Solids < 10.0%	Use professional judgment	Use professional judgment
10.0% ≤ %Solids ≤ 30.0%	Use professional judgment	Use professional judgment
%Solids > 30.0%	No qualification	No qualification

SAMPLE QUANTITATION

The sample quantitation evaluation is to verify laboratory quantitation results. In the space below, please show a minimum of one sample calculation:

Sample ID:__ FA42192-3_____ Analyte:__1,4-dioxane__ RF:_0.599_

[ ] = (80516)(4.0)/(28666)(0.599)
= 18.8 ppm Ok

DATA REVIEW WORKSHEETS

QUANTITATION LIMITS

A. Dilution performed

SAMPLE ID	DILUTION FACTOR	REASON FOR DILUTION
FA42192-12	10 x	1,4-Dioxane over calibration range
FA42192-13	25 x	1,4-Dioxane over calibration range
FA42192-6	10 x	Matrix interference

All criteria were met __X__
Criteria were not met
and/or see below _____

FIELD DUPLICATE PRECISION

Sample IDs:_FA42192-8/FA42192-11_____ Matrix:__Groundwater_____

Field duplicates samples may be taken and analyzed as an indication of overall precision. These analyses measure both field and lab precision; therefore, the results may have more variability than laboratory duplicates which only laboratory performance. It is also expected that soil duplicate results will have a greater variance than water matrices due to difficulties associated with collecting identical field duplicate samples.

The project QAPP should be reviewed for project-specific information.
Suggested criteria: if large RPD (> 50 %) is observed, confirm identification of the samples and note differences. If both samples and duplicate are <5 SQL, the RPD criteria is doubled.

COMPOUND	SQL ug/L	SAMPLE CONC. (ug/l)	DUPLICATE CONC. (ug/l)	RPD	ACTION
Field duplicate analyzed as part of this data package. RPD within the required guidance document criteria < 50 % for detected target analytes above 5 SQL.					

All criteria were met __X__
Criteria were not met
and/or see below _____

OTHER ISSUES

A. System Performance

List samples qualified based on the degradation of system performance during simple analysis:

Sample ID	Comments	Actions

Action:

Use professional judgment to qualify the data if it is determined that system performance has degraded during sample analyses. Inform the Contract Laboratory Program COR any action as a result of degradation of system performance which significantly affected the data.

B. Overall Assessment of Data

List samples qualified based on other issues:

Sample ID	Comments	Actions

_No_other_issues_that_required_the_need_to_qualify_the_data._Results_are_valid_and_can_be_used_for_decission_purposes._Other_discrepancies_are_shown_below._____

Note:

Action:

1. Use professional judgment to determine if there is any need to qualify data which were not qualified based on the Quality Control (QC) criteria previously discussed.
2. Write a brief narrative to give the user an indication of the analytical limitations of the data. Inform the Contract Laboratory COR the action, any inconsistency of the data with the Sample Delivery Group (SDG) Narrative. If sufficient information on the intended use and required quality of the data is available, the reviewer should include their assessment of the usability of the data within the given context. This may be used as part of a formal Data Quality Assessment (DQA).

3. Sometimes, due to dilutions, re-analysis or SIM/Scan runs are being performed, there will be multiple results for a single analyte from a single sample. The following criteria and professional judgment are used to determine which result should be reported:
    - The analysis with the lower CRQL
    - The analysis with the better QC results
    - The analysis with the higher results

EXECUTIVE NARRATIVE

SDG No:	**FA42192**	Laboratory:	**Accutest, Orlando**
Analysis:	**MADEP VPH**	Number of Samples:	**11**
Location:	BMSMC, Humacao, PR		

SUMMARY: Eleven (11) samples were analyzed for Volatiles TPHC Ranges by method MADEP VPH. Samples were validated following the METHOD FOR THE DETERMINATION OF VOLATILE PETROLEUM HYDROCARBONS (VPH) quality control criteria, Massachusetts Department of Environmental Protection, Revision 1.1 (2004). Also the general validation guidelines promulgated by the USEPA Hazardous Wastes Support Section. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

Results are valid and can be used for decision making purposes.

Critical issues: **None**
Major: **None**
Minor: **None**

Critical findings: **None**
Major findings: **None**
Minor findings: **1.** As per chain-of-custody form sample FA42192-3 was to use as the QC sample (MS/MSD) for this batch. Samples FA42192-2 and FA42237-4 were used instead. No action taken, professional judgment.

MS/MSD % recovery and RPD within laboratory control limits except for the cases described in this document. No action taken, results apply to unspiked sample. Unspiked sample was from another batch.

COMMENTS: Results are valid and can be used for decision making purposes.

Reviewers Name: Rafael Infante
Chemist License 1888

Signature: Rafael Infante

Date: April 26, 2017

ORGANIC DATA SAMPLE SUMMARY

Sample ID: FA42192-1
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: AQ - Field Blank Water

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	100	ug/L	1.0	-	U	Yes

Sample ID: FA42192-2
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	100	ug/L	1.0	-	U	Yes

Sample ID: FA42192-3
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	35.6	ug/L	1.0	J	J	Yes

Sample ID: FA42192-6
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	38.4	ug/L	1.0	J	J	Yes
C9 - C12 Aliphatics (Unadj.)	74500	ug/L	500	-	-	Yes
C9 - C10 Aromatics (Unadj.)	262	ug/L	1.0	-	-	Yes

Sample ID: FA42192-7
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	100	ug/L	1.0	-	U	Yes

Sample ID: FA42192-8
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	100	ug/L	1.0	-	U	Yes

Sample ID: FA42192-11
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	100	ug/L	1.0	-	U	Yes

Sample ID: FA42192-12
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	100	ug/L	1.0	-	U	Yes

Sample ID: FA42192-13
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	100	ug/L	1.0	-	U	Yes

Sample ID: FA42192-14
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	100	ug/L	1.0	-	U	Yes

Sample ID: FA42192-15
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: AQ - Field Blank Water

METHOD: MADEP VPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C5 - C8 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C12 Aliphatics (Unadj.)	100	ug/L	1.0	-	U	Yes
C9 - C10 Aromatics (Unadj.)	100	ug/L	1.0	-	U	Yes

DATA REVIEW WORKSHEETS

Type of validation Full:__X____ Limited:_____

Project Number:_FA42192_____
Date:_____03/16-17/2017_____
Shipping date:_____03/17/2017_____
EPA Region:_____2_____

REVIEW OF VOLATILE PETROLEUM HYDROCARBON (VPHs) PACKAGE

The following guidelines for evaluating volatile organics were created to delineate required validation actions. This document will assist the reviewer in using professional judgment to make more informed decision and in better serving the needs of the data users. The sample results were assessed according to the data validation guidance documents in the following order of precedence METHOD FOR THE DETERMINATION OF VOLATILE PETROLEUM HYDROCARBONS (VPH), Massachusetts Department of Environmental Protection, Revision 1.1 (2004). Also the general validation guidelines promulgated by the USEPA Hazardous Wastes Support Section. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

The hardcopied (laboratory name) _Accutest_Laboratories_-_Orlando_____ data package received has been reviewed and the quality control and performance data summarized. The data review for VOCs included:

Lab. Project/SDG No.: __FA42192_____ Sample matrix: ____Groundwater_____
No. of Samples: _____11_____
Field blank No.: _____ FA42192-1;_FA42192-15_____
Equipment blank No.: ___-_____
Trip blank No.: _____-_____
Field duplicate No.: _____-_____

__X____ Data Completeness
__X____ Holding Times
__N/A__ GC/MS Tuning
__N/A__ Internal Standard Performance
__X____ Blanks
__X____ Surrogate Recoveries
__X____ Matrix Spike/Matrix Spike Duplicate
___X__ Laboratory Control Spikes
___X__ Field Duplicates
___X__ Calibrations
___X__ Compound Identifications
___X__ Compound Quantitation
___X__ Quantitation Limits

Overall Comments: _Volatiles_by_GC_by_Method_MADEP_VPH,_REV_1.1._(C9- C10 Aromatics (Unadj.)).

Definition of Qualifiers:

J- Estimated results
U- Compound not detected
R- Rejected data
UJ- Estimated nondetect

Reviewer: Rafael Infante
Date:_____April_24,_2017_____

All criteria were met __x___
Criteria were not met and/or see below _____

I. DATA COMPLETNESS
A. Data Package:

MISSING INFORMATION	DATE LAB. CONTACTED	DATE RECEIVED

B. Other Discrepancies:

All criteria were met ___X___
Criteria were not met and/or see below _____

HOLDING TIMES
The objective of this parameter is to ascertain the validity of the results based on the holding time of the sample from time of collection to the time of extraction, and subsequently from the time of extraction to the time of analysis.

Complete table for all samples and note the analysis and/or preservation not within criteria

SAMPLE ID	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	ACTION
Samples analyzed within method recommended holding time. Sample preservation within the required criteria.				

Criteria

Preservation:

Samples analyzed with ambient purge temperature: Samples must be acidified to a pH of 2.0 or less at the time of collection.
Samples analyzed with heated purge temperature: Samples must be treated to a pH of 11.0 or greater at the time of collection.
Methanol preservation of soil/sediment samples is mandatory. Methanol (purge-and-trap grade) must be added to the sample vial before or immediately after sample collection. In lieu of the in-field preservation of samples with methanol, soil samples may be obtained in specially-designed air tight sampling devices, provided that the samples are extruded and preserved in methanol within 48 hours of collection.

Holding times:

Aqueous samples using ambient or heated purge - analyze within 14 days.
Soil/sediment samples - analysis within 28 days.

Cooler temperature (Criteria: 4 ± 2 °C):_____1.6-5.8_°C_____

Actions: Qualify positive results/non-detects as follows:

If holding times are exceeded, estimate positive results (J) and nondetects (UJ).
If holding times are grossly exceeded, use professional judgment to qualify data. The data reviewer may choose to estimate positive results (J) and rejects nondetects (R).
If samples were not at the proper temperature (> 10°C) or improperly preserved, use professional judgment to qualify the results.

All criteria were met ___X___
Criteria were not met and/or see below _____

CALIBRATIONS VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:___03/21/17_____

Dates of initial calibration verification:__03/21/17_____

Instrument ID numbers:_____HP5890_____

Matrix/Level:_____AQUEOUS/MEDIUM_____

DATE	LAB FILE ID#	ANALYTE	CRITERIA OUT RFs, %RSD, %D, r	SAMPLES AFFECTED
Initial and initial calibration verification meet method specific requirements				

Criteria- ICAL

- Five point calibration curve.
- The percent relative standard deviation (%RSD) of the calibration factor must be equal to or less than 25% over the working range for the analyte of interest. When this condition is met, linearity through the origin may be assumed, and the average calibration factor is used in lieu of a calibration curve.
- A collective calibration factor must also be established for each hydrocarbon range of interest. Calculate the collective CFs for C5-C8 Aliphatic Hydrocarbons and C9-C12 Aliphatic Hydrocarbons using the FID chromatogram. Calculate the collective CF for the C9-C10 Aromatic Hydrocarbons using the PID chromatogram. Tabulate the summation of the peak areas of all components in that fraction against the total concentration injected. The %RSD of the calibration factor must be equal to or less than 25% over the working range for the hydrocarbon range of interest.

Criteria- CCAL

- At a minimum, the working calibration factor must be verified on each working day, after every 20 samples, and at the end of the analytical sequence by the injection of a mid-level continuing calibration standard to verify instrument performance and linearity.
- If the percent difference (%D) for any analyte varies from the predicted response by more than ±25%, a new five-point calibration must be performed for that analyte. Greater percent differences are permissible for n-nonane. If the %D for n-nonane is greater than 30, note the nonconformance in the case narrative. It should be noted that the %Ds are calculated when CFs are used for the initial calibration and

percent drifts are calculated when calibration curves using linear regression are used for the initial calibration.

Actions:

If %RSD > 25% for target compounds or a correlation coefficient < 0.99, estimate positive results (J) and use professional judgment to qualify nondetects.
If % D > 25% (> 30 for nonane), estimate positive results (J) and nondetects (UJ).

CALIBRATIONS VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:_____03/21/17_____
Dates of continuing calibration verification:__03/26/17;_03/27/17_____
Dates of final calibration verification:_____03/27/17:_03/27/17_____
Instrument ID numbers:_____VOA10_____
Matrix/Level:_____AQUEOUS/MEDIUM_____

DATE	LAB FILE ID#	ANALYTE	CRITERIA OUT RFs, %RSD, %D, r	SAMPLES AFFECTED

Note: Continuing and final calibration verification meets method specific requirements.

A separate worksheet should be filled for each initial curve

All criteria were met ___X___
Criteria were not met and/or see below _____

V A. BLANK ANALYSIS RESULTS (Sections 1 & 2)

The assessment of the blank analysis results is to determine the existence and magnitude of contamination problems. The criteria for evaluation of blanks apply only to blanks associated with the samples, including trip, equipment, and laboratory blanks. If problems with any blanks exist, all data associated with the case must be carefully evaluated to determine whether or not there is an inherent variability in the data for the case, or if the problem is an isolated occurrence not affecting other data. A Laboratory Method Blank must be run after samples suspected of being highly contaminated to determine if sample carryover has occurred.

List the contamination in the blanks below. High and low levels blanks must be treated separately.

Laboratory blanks

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS

_METHOD_BLANKS_MEET_THE_METHOD_SPECIFIC_CRITERIA_____
_____
_____
_____
_____
_____

Note:

Field/Trip/Equipment

A methanol trip blank or acidified reagent water trip blank **should** continually accompany each soil/sediment sample or water sample batch, respectively, during sampling, storage, and analysis.

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS

_NO_TRIP/EQUIPMENT_BLANKS_ASSOCIATED_WITH_THIS_DATA_PACKAGE._____
_NO_TARGET_ANALYTE_DETECTED_IN_THE_FIELD_BLANKS_ASSOCIATED_WITH_
_THIS_DATA_PACKAGE._____
_____
_____
_____

Note:

V B. BLANK ANALYSIS RESULTS (Section 3)

Blank Actions

The ALs for samples which have been diluted should be corrected for the sample dilution factor and/or % moisture, where applicable. Peaks must not be detected above the Reporting Limit within the retention time window of any analyte of interest. The hydrocarbon ranges must not be detected at a concentration greater than 10% of the most stringent MCP cleanup standard. Specific actions area as follows:

If the concentration is < sample quantitation limit (SQL) and < AL, report the compound as not detected (U) at the SQL.
If the concentration is ≥ SQL but < AL, report the compound as not detected (U) at the reported concentration.
If the concentration is > AL, report the concentration unqualified.

All criteria were met __X___
Criteria were not met and/or see below _____

SURROGATE SPIKE RECOVERIES

Laboratory performance of individual samples is established by evaluation of surrogate spike recoveries. All samples are spiked with surrogate compounds prior to sample analysis. The accuracy of the analysis is measured by the surrogate percent recovery. Since the effects of the sample matrix are frequently outside the control of the laboratory and may present relatively unique problems, the validation of data is frequently subjective and demands analytical experience and professional judgment.
List the percent recoveries (%Rs) which do not meet the criteria for surrogate recovery.
Matrix: solid/aqueous

SAMPLE ID	SURROGATE COMPOUND	ACTION
	BFB	

_____
_SURROGATE_STANDARD_RECOVERIES_WITHIN_LABORATORY_CONTROL_____
_LIMITS._____
_____
_____
_____
_____

QC Limits* (Aqueous)
_____LL_to_UL___ _70_to_130_ ___to_____ ___to_____
QC Limits* (Solid)
_____LL_to_UL___ _70_to_130_ ___to_____ ___to_____

It is recommended that surrogate standard recoveries be monitored and documented on a continuing basis. At a minimum, when surrogate recovery from a sample, blank, or QC sample is less than 70% or more than 130%, check calculations to locate possible errors, check the fortifying standard solution for degradation, and check changes in instrument performance.

If the cause cannot be determined, reanalyze the sample unless one of the following exceptions applies:

(1) Obvious interference is present on the chromatogram (e.g., unresolved complex mixture);
(2) Percent moisture of associated soil/sediment sample is >25% and surrogate recovery is >10%; or
(3) The surrogate exhibits high recovery and associated target analytes or hydrocarbon ranges are not detected in sample.

If a sample with a surrogate recovery outside of the acceptable range is not reanalyzed based on any of these aforementioned exceptions, this information must be noted on the data report form and discussed in the Executive Report. Analysis of the sample on dilution may diminish matrix-related surrogate recovery problems. This approach can be used as long as the reporting limits to evaluate applicable MCP standards can still be achieved with the dilution. If not, reanalysis without dilution must be performed.

All criteria were met _____
Criteria were not met and/or see below __X___

VII. A MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

This data is generated to determine long term precision and accuracy in the analytical method for various matrices. This data alone cannot be used to evaluate the precision and accuracy of individual samples.

At the request of the data user, and in consideration of sample matrices and data quality objectives, matrix spikes and matrix duplicates may be analyzed with every batch of 20 samples or less per matrix.

- **Matrix duplicate** - Matrix duplicates are prepared by analyzing one sample in duplicate. The purpose of the matrix duplicates is to determine the homogeneity of the sample matrix as well as analytical precision. The RPD of detected results in the matrix duplicate samples must not exceed 50 when the results are greater than 5x the reporting limit.
- The desired spiking level is 50% of the highest calibration standard. However, the total concentration in the MS (including the MS and native concentration in the unspiked sample) should not exceed 75% of the highest calibration standard in order for a proper evaluation to be performed. The purpose of the matrix spike is to determine whether the sample matrix contributes bias to the analytical results. The corrected concentrations of each analyte within the matrix spiking solution must be within 70 - 130% of the true value. Lower recoveries of n-nonane are permissible (if included in the calibration of the C9-C12 aliphatic range), but must be noted in the narrative if <30%.

MS/MSD Recoveries and Precision Criteria

Sample ID:_FA42237-4_MS/MSD_____ Matrix/Level:_Groundwater____

List the %Rs, RPD of the compounds which do not meet the QC criteria.

The QC reported here applies to the following samples:
FA42192-1, FA42192-2, FA42192-3, FA42192-6, FA42192-7, FA42192-8, FA42192-11, FA42192-12, FA42192-13, FA42192-14, FA42192-15

Compound	FA42237-4 ug/l	Q	Spike ug/l	MS ug/l	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
C5- C8 Aliphatics (Unadj.)	ND		480	262	55*	480	262	55*	0	70-130/50
C9- C12 Aliphatics (Unadj.)	ND		400	151	38*	400	153	38*	1	70-130/50
C9- C10 Aromatics (Unadj.)	ND		240	73.8	31*	240	82.6	34*	11	70-130/50

* Outside laboratory control limits.

Note: As per chain-of-custody form sample FA42192-3 was to use as the QC sample (MS/MSD) for this batch. Samples FA42192-2 and FA42237-4 were used instead. No action taken, professional judgment.

MS/MSD % recovery and RPD within laboratory control limits except for the cases described in this document. No action taken, results apply to unspiked sample. Unspiked sample was from another batch.

No action is taken on MS/MSD results alone to qualify the entire case. However, used informed professional judgment, the data reviewer may use the MS/MSD results in conjunction with other QC criteria and determine the need for some qualification of the data. In those instances where it can be determined that the results of the MS/MSD affect only the sample spiked, the qualification should be limited to this sample alone. However, it may be determined through the MS/MSD results that the laboratory is having a systematic problem in the analysis of one or more analytes, which affects the associated samples.

2. MS/MSD – Unspiked Compounds

List the concentrations of the unspiked compounds and determine the % RSDs of these compounds in the unspiked sample, matrix spike, and matrix spike duplicate.

COMPOUND	CONCENTRATION SAMPLE	MS	MSD	%RPD	ACTION

Criteria: None specified, use %RSD $\leq$ 50 as professional judgment.

Actions:

If the % RSD > 50, qualify the results in the spiked sample as estimate (J).
If the % RSD is not calculable (NC) due to nondetect value in the sample, MS, and/or MSD, use professional judgment to qualify sample data.

A separate worksheet should be used for each MS/MSD pair.

All criteria were met __X___
Criteria were not met and/or see below _____

VIII. LABORATORY CONTROL SAMPLE (LCS/LCSD) ANALYSIS

This data is generated to determine accuracy of the analytical method for various matrices.

1. LCS Recoveries Criteria

List the %R of compounds which do not meet the criteria

LCS ID	COMPOUND	% R	QC LIMIT	ACTION

__LCS/LCSD_RECOVERY_WITHIN_LABORATORY_CONTROL_LIMTS._____
_____
_____
_____
_____
_____

Criteria:
* Refer to QAPP for specific criteria.
* The spike recovery must be between 70% and 130%. Lower recoveries of n-nonane are permissible (if included in the calibration of the C9-C12 aliphatic range). If the recovery of n-nonane is <30%, note the nonconformance in the executive narrative.

Actions:
Actions on LCS recovery should be based on both the number of compounds that are outside the %R criteria and the magnitude of the excedance of the criteria.

If the %R of the analyte is > UL, qualify all positive results (j) for the affected analyte in the associated samples and accept nondetects.
If the %R of the analyte is < LL, qualify all positive results (j) and reject (R) nondetects for the affected analyte in the associated samples.
If more than half the compounds in the LCS are not within the required recovery criteria, qualify all positive results as (J) and reject nondetects (R) for all target analyte(s) in the associated samples.

2. Frequency Criteria:

Where LCS analyzed at the required frequency and for each matrix (1 per 20 samples per matrix)? **Yes** or No.
If no, the data may be affected. Use professional judgment to determine the severity of the effect and qualify data accordingly. Discuss any actions below and list the samples affected. Discuss the actions below:

All criteria were met __X___
Criteria were not met and/or see below _____

IX. FIELD/LABORATORY DUPLICATE PRECISION

Sample IDs:__FA42196-6/-6DUP_____ Matrix:__Groundwater__
Sample IDs:__FA42196-8/FA42196-11____ Matrix:__Groundwater__

Field/laboratory duplicates samples may be taken and analyzed as an indication of overall precision. These analyses measure both field and lab precision; therefore, the results may have more variability than laboratory duplicates which measures only laboratory performance. It is also expected that soil duplicate results will have a greater variance than water matrices due to difficulties associated with collecting identical field duplicate samples.

COMPOUND	SQL	SAMPLE CONC.	DUPLICATE CONC.	RPD	ACTION
Field/laboratory duplicate analyzed with this data package. RPD within laboratory and validation guidance document criteria (± 50 %) for analytes detected above reporting limits.					

Criteria:

The project QAPP should be reviewed for project-specific information.
RPD ± 30% for aqueous samples, RPD ± 50 % for solid samples if results are ≥ SQL.
If both samples and duplicate are <5 SQL, the RPD criteria is doubled.

SQL = soil quantitation limit

Actions:

If both the sample and the duplicate results are nondetects (ND), the RPD is not calculable (NC). No action is needed.

Qualify as estimated positive results (J) and nondetects (UJ) for the compound that exceeded the above criteria.

If one sample result is not detected and the other is ≥ 5x the SQL qualify (J/UJ).

Note: If SQLs for the sample and duplicate are significantly different, use professional judgment to determine if qualification is appropriate.

If one sample value is not detected and the other is < 5x the SQL, use professional judgment to determine if qualification is appropriate.

All criteria were met __X___
Criteria were not met and/or see below _____

XI. COMPOUND IDENTIFICATION

The compound identification evaluation is to verify that the laboratory correctly identified target analytes as well as tentatively identified compounds (TICs).

1. Verify that the target analytes were within the retention time windows.

   - Retention time windows must be re-established for each Target VPH Analyte each time a new GC column is installed, and must be verified and/or adjusted on a daily basis.
   - Coelution of the m- and p- xylene isomers is permissible.
   - All surrogates must be adequately resolved from individual Target Analytes included in the VPH Component Standard.
   - For the purposes of this method, adequate resolution is assumed to be achieved if the height of the valley between two peaks is less than 25% of the average height of the two peaks.
   - The n-pentane (C5) and MtBE peaks must be adequately resolved from any solvent front that may be present on the FID and PID chromatograms, respectively.

   Note: Target analytes were within the retention time window.

2. If target analytes and/or TICs were not correctly identified, request that the laboratory resubmit the corrected data.

All criteria were met __X___
Criteria were not met and/or see below _____

XII. QUANTITATION LIMITS AND SAMPLE RESULTS

The sample quantitation evaluation is to verify laboratory quantitation results.

1. In the space below, please show a minimum of one sample calculation:

FID

Computer printout

PID

Computer printout

2. If requested, verify that the results were above the laboratory method detection limit (MDLs).

3. If dilutions performed, were the SQLs elevated accordingly by the laboratory? List the affected samples and dilution factor in the table below.

SAMPLE ID	DILUTION FACTOR	REASON FOR DILUTION
FA42196-6	500 x	C9 – C12 aliphatics (unadj.) above calibration concentration range.

If dilution was not performed and the results were above the concentration range, estimate results (J) for the affected compounds. List the affected samples/compounds:

_____

_____

EXECUTIVE NARRATIVE

SDG No:	**FA42192**	Laboratory:	**Accutest, New Jersey**
Analysis:	**SW846-8015C**	Number of Samples:	**11**

Location: BMSMC, Humacao, PR

SUMMARY: Eleven (11) samples were analyzed for the low molecular weight alcohols (LMWAs) list following method SW846-8015C. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence: "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846 (Final Update III, December 1996)," specifically for Methods 8000/8015C are utilized. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

Results are valid and can be used for decision making purposes.

Critical issues: **None**
Major: **None**
Minor: **None**

Critical findings: **None**
Major findings: **None**
Minor findings: **1.** Isopropyl alcohol and sec-butyl alcohol recovered high in laboratory control sample (Blank Spike). No action taken, professional judgment. High percent recoveries and no associated positive reported in the QC batch.

COMMENTS: Results are valid and can be used for decision making purposes.

Reviewers Name: Rafael Infante
Chemist License 1888

Signature: Rafael Infante
Date: April 26, 2017

SAMPLE ORGANIC DATA SAMPLE SUMMARY - JC18972

Sample ID: FA42192-1
Sample location: BMSM, Humacao, PR
Sampling date: 3/16/2017
Matrix: AQ - Field Blank Water

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-2
Sample location: BMSM, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-3
Sample location: BMSM, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-6
Sample location: BMSM, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-7
Sample location: BMSM, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-8
Sample location: BMSM, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-11
Sample location: BMSM, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-12
Sample location: BMSM, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-13
Sample location: BMSM, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-14
Sample location: BMSM, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

Sample ID: FA42192-15
Sample location: BMSM, Humacao, PR
Sampling date: 3/17/2017
Matrix: AQ - Field Blank Water

METHOD: 8015C

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Ethanol	100	ug/l	1.0	-	U	Yes
Isobutyl Alcohol	100	ug/l	1.0	-	U	Yes
Isopropyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Propyl Alcohol	100	ug/l	1.0	-	U	Yes
n-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
sec-Butyl Alcohol	100	ug/l	1.0	-	U	Yes
Methanol	200	ug/l	1.0	-	U	Yes

DATA REVIEW WORKSHEETS

Project Number:___FA42192_____
Date:_____03/16-17/2017_____
Shipping Date:___03/17/2017_____
EPA Region:_____2_____

REVIEW OF VOLATILE ORGANIC PACKAGE

The following guidelines for evaluating volatile organics were created to delineate required validation actions. This document will assist the reviewer in using professional judgment to make more informed decision and in better serving the needs of the data users. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence: "Test Methods for Evaluating Solid Waste, Physical/Chemical Methods SW-846 (Final Update III, December 1996)," specifically for Methods 8000/8015C are utilized. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

The hardcopied (laboratory name) _Accutest,_New_Jersey_____ data package received has been reviewed and the quality control and performance data summarized. The modified data review for VOCs included:

Lab. Project/SDG No.: __FA42192_____ Sample matrix: ___Groundwater_____
No. of Samples: _____11_____

Trip blank No.: _____-_____
Field blank No.: _____FA42192-11;_FA42192-15_____
Equipment blank No.:_____
Field duplicate No.:_____FA42192-8/FA42192-11_____

___X___ Data Completeness
___X___ Holding Times
___N/A_ GC/MS Tuning
___N/A_ Internal Standard Performance
___X___ Blanks
___X___ Surrogate Recoveries
___X___ Matrix Spike/Matrix Spike Duplicate
___X___ Laboratory Control Spikes
___X___ Field Duplicates
___X___ Calibrations
___X___ Compound Identifications
___X___ Compound Quantitation
___X___ Quantitation Limits

Overall Comments:_Low_molecular_weight_alcohols_by_SW-846_8015C._____
_____
_____

Definition of Qualifiers:

J- Estimated results
U- Compound not detected
R- Rejected data
UJ- Estimated nondetect

Reviewer:_____
Date:__April_24,_2017_____

DATA REVIEW WORKSHEETS

DATA COMPLETENESS

MISSING INFORMATION	DATE LAB. CONTACTED	DATE RECEIVED

All criteria were met __X__
Criteria were not met
and/or see below _____

HOLDING TIMES

The objective of this parameter is to ascertain the validity of the results based on the holding time of the sample from time of collection to the time of analysis.

Complete table for all samples and note the analysis and/or preservation not within criteria

SAMPLE ID	DATE SAMPLED	DATE ANALYZED	pH	ACTION
All samples analyzed within the recommended method holding. All samples properly preserved.				

Criteria

Aqueous samples – 14 days from sample collection for preserved samples (pH $\leq$ 2, 4°C), no air bubbles.
Aqueous samples – 7 days from sample collection for unpreserved samples, 4°C, no air bubbles.
Groundwater samples- 7 days from sample collection.
Cooler temperature (Criteria: 4 $\pm$ 2 °C): 1.6 – 5.8°C

Actions

If the VOCs vial(s) have air bubbles, estimate positive results (J) and reject nondetects (R).
If the % solids of Groundwater samples is 10-50%, estimates positive results (J) and nondetects (UJ)
If the % solid of Groundwater samples is < 10%, estimate positive results (J) and reject nondetects (R).
If holding times are exceeded but < 14 days beyond criteria, estimate positive results (J) and nondetects (UJ).
If holding times are exceeded but < 28 days beyond criteria, estimate positive results (J) and reject nondetects (R).
If holding times are grossly exceeded (> 28 days beyond criteria), reject all results (R).
If samples were not iced or if the ice were melted (> 10°C), estimate positive results (J) and nondetects (UJ).

All criteria were met __N/A__
Criteria were not met see below ____

GC/MS TUNING

The assessment of the tuning results is to determine if the sample instrumentation is within the standard tuning QC limits

__N/A_ The BFB performance results were reviewed and found to be within the specified criteria.

__N/A_ BFB tuning was performed for every 12 hours of sample analysis.

If no, use professional judgment to determine whether the associated data should be accepted, qualified or rejected.

List the samples affected: _____

If mass calibration is in error, all associated data are rejected.

DATA REVIEW WORKSHEETS

All criteria were met __X__
Criteria were not met
and/or see below _____

CALIBRATION VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:_____10/10/16_____
Dates of continuing calibration:_____03/24/17;_03/27/17_____
Dates of final calibration verification:___03/24/17;_03/27/17_____
Instrument ID number:_____GCGH_____
Matrix/Level:_____Aqueous/low_____

DATE	LAB FILE ID#	CRITERIA OUT RFs, %RSD, %D, r	COMPOUND	SAMPLES AFFECTED

Note: Initial, continuing, and final calibration verifications meets method specific criteria in at least one of the two columns. Final calibration verification included in data packages.

Criteria

All RFs must be > 0.05 regardless of method requirements for SPCC.
All %RSD must be ≤ 15 % regardless of method requirements for CCC.
All %Ds must be ≤ 20% regardless of method requirements for CCC.
It should be noted that Region 2 SOP HW-24 does not specify criterion for the curve correlation coefficient (r). A limit for r of ≥ 0.995 has therefore been utilized as professional judgment.

Actions

If any compound has an initial RF or a continuing RF of < 0.05, estimate positive results (J) and reject nondetects (R), regardless of method requirements.
If any compound has a %RSD > 15%, estimate positive results (J) and use professional judgment to qualify nondetects.
If any compound has a %RSD > 90%, estimate positive results (J) and reject nondetects (R).
If any compound has a % D > 20%, estimate positive results (J) and reject nondetects (R).
If any compound has a % D > 20%, estimate positive results (J) and nondetects (UJ).
If any compound has a % D > 90%, estimate positive results (J) and reject nondetects (R).
If any compound has r < 0.995, estimate positive results and nondetects.

A separate worksheet should be filled for each initial curve

All criteria were met __X__
Criteria were not met
and/or see below _____

V A. BLANK ANALYSIS RESULTS (Sections 1 & 2)

The assessment of the blank analysis results is to determine the existence and magnitude of contamination problems. The criteria for evaluation of blanks apply only to blanks associated with the samples, including trip, equipment, and laboratory blanks. If problems with any blanks exist, all data associated with the case must be carefully evaluated to determine whether or not there is an inherent variability in the data for the case, or if the problem is an isolated occurrence not affecting other data.

List the contamination in the blanks below. High and low levels blanks must be treated separately.

Laboratory blanks

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
____All_method_blank_meeth_method_specific_criteria_____				

Field/Equipment/Trip blank

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS

_____
_No_target_analytes_detected_in_the_field_blanks_analyzed_with_this_data_package._No_equipment/
_trip_blanks_analyzed_with_this_data_package._____

All criteria were met __X__
Criteria were not met
and/or see below _____

V B. BLANK ANALYSIS RESULTS (Section 3)

Blank Actions

Action Levels (ALs) should be based upon the highest concentration of contaminant determined in any blank. Do not qualify any blank with another blank. The ALs for samples which have been diluted should be corrected for the sample dilution factor and/or % moisture, where applicable. No positive sample results should be reported unless the concentration of the compound in the samples exceeds the ALs:

ALs = 10x the amount of common contaminants (methylene chloride, acetone, 2-butanone, and toluene)
ALs = 5x for any other compounds

Specific actions are as follows:

If the concentration is < sample quantitation limit (SQL) and ≤ AL, report the compound as not detected (U) at the SQL.
If the concentration is ≥ SQL but ≤ AL, report the compound as not detected (U) at the reported concentration.
If the concentration is ≥ SQL and > AL, report the concentration unqualified.

Notes:

High and low level blanks must be treated separately
Compounds qualified "U" for blank contamination are still considered "hits" when qualifying for calibration criteria.

CONTAMINATION SOURCE/LEVEL	COMPOUND	CONC/UNITS	AL/UNITS	SQL	AFFECTED SAMPLES

All criteria were met ___X___
Criteria were not met
and/or see below _____

SURROGATE SPIKE RECOVERIES

Laboratory performance of individual samples is established by evaluation of surrogate spike recoveries. All samples are spiked with surrogate compounds prior to sample analysis. The accuracy of the analysis is measured by the surrogate percent recovery. Since the effects of the sample matrix are frequently outside the control of the laboratory and may present relatively unique problems, the validation of data is frequently subjective and demands analytical experience and professional judgment.
List the percent recoveries (%Rs) which do not meet the criteria for surrogate recovery.
Matrix: solid/aqueous

SAMPLE ID	**SURROGATE COMPOUND**				**ACTION**
	Hexanol	**~~DBFM~~**	**~~TOL-d8~~**	**~~BFB~~**	
	S1 a	S1 b			
FA42192-1	79	78			
FA42192-2	69	68			
FA42192-3	79	73			
FA42192-6	78	75			
FA42192-7	81	76			
FA42192-8	80	78			
FA42192-11	77	76			
FA42192-12	75	75			
FA42192-13	72	77			
FA42192-14	105	80			
FA42192-15	75	72			
GGH5693-BS	99	104			
GGH5693-MB2	93	91			
GGH5695-BS	111	111			
GGH5695-MB1	87	83			
JC38823-4MS	85	68			
JC38823-4MSD	80	73			
JC38825-3MS	85	86			
JC38825-3MSD	88	84			
GGH5693-MB1	105	112			

(a) Recovery from GC signal #2 (b) Recovery from GC signal #1

Note: All surrogate recoveries within laboratory control limits.

QC Limits* (Aqueous)
_____LL_to_UL___ _56_to_145_ ___to_____ ___to_____ ___to_____
QC Limits* (Solid-Low)
_____LL_to_UL___ ____to_____ ___to_____ ___to_____ ___to_____
QC Limits* (Solid-Med)
_____LL_to_UL___ ____to_____ ___to_____ ___to_____ ___to_____

* QC limits are laboratory in-house performance criteria, LL = lower limit, UL = upper limit.
* If QC limits are not available, use limits of 80 – 120 % for aqueous and 70 – 130 % for solid samples.

Actions:

QUALITY	%R < 10%	%R = 10% - LL	%R > UL
Positive results	J	J	J
Nondetects results	R	UJ	Accept

Surrogate action should be applied:

If one or more surrogate in the VOC fraction is out of specification, but has a recovery of > 10%.
If any one surrogate in a fraction shows < 10 % recovery.

All criteria were met __X____
Criteria were not met
and/or see below _____

VII. A MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

This data is generated to determine long term precision and accuracy in the analytical method for various matrices. This data alone cannot be used to evaluate the precision and accuracy of individual samples. If any % R in the MS or MSD falls outside the designated range, the reviewer should determine if there are matrix effects, i.e. LCS data are within the QC limits but MS/MSD data are outside QC limit.

1. MS/MSD Recoveries and Precision Criteria

The laboratory should use one MS and a duplicate analysis of an unspiked field sample if target analytes are expected in the sample. If target analytes are not expected, MS/MSD should be analyzed.

List the %Rs, RPD of the compounds which do not meet the criteria.

Sample ID:__JC38825-3MS/-3MSD_____ Matrix/Level:____Groundwater/medium_____
Sample ID:__JC38823-4MS/-4MSD_____ Matrix/Level:____Groundwater/medium_____

MS OR MSD	COMPOUND	% R	RPD	QC LIMITS	ACTION

MS/MSD%_recoveries_and_RPD_within_laboratory_control_limits._____

Note:

* QC limits are laboratory in-house performance criteria, LL = lower limit, UL = upper limit.
* If QC limits are not available, use limits of 70 – 130 %.

Actions:

QUALITY	%R < LL	%R > UL
Positive results	J	J
Nondetects results	R	Accept

All criteria were met __X__
Criteria were not met
and/or see below _____

MS/MSD criteria apply only to the unspiked sample, its dilutions, and the associated MS/MSD samples:

If the % R for the affected compounds were < LL (or 70 %), qualify positive results (J) and nondetects (UJ).
If the % R for the affected compounds were > UL (or 130 %), only qualify positive results (J).
If 25 % or more of all MS/MSD %R were < LL (or 70 %) or if two or more MS/MSD %Rs were < 10%, qualify all positive results (J) and reject nondetects (R).

VII. B MATRIX SPIKE/MATRIX SPIKE DUPLICATE

MS/MSD – Unspiked Compounds

It should be noted that Region 2 SOP HW-24 does not specify a MS/MSD criteria for the unspiked compounds in the sample. A %RSD of < 50% has therefore been utilized as professional judgment.

If all target analytes were spiked in the MS/MSD, this review element is not applicable.

List the %RSD of the compounds which do not meet the criteria.

Sample ID:_____-_____ Matrix/Level/Unit:_____-_____

COMPOUND	SAMPLE CONC.	MS CONC.	MSD CONC.	% RSD	ACTION

Actions:

* If the % RSD > 50, qualify the positive result in the unspiked samples as estimated (J).
* If the % RSD is not calculated (NC) due to nondetected value, use professional judgment to qualify the data.

A separate worksheet should be used for each MS/MSD pair.

All criteria were met __X__
Criteria were not met
and/or see below _____

VIII. LABORATORY CONTROL SAMPLE (LCS) ANALYSIS

This data is generated to determine accuracy of the analytical method for various matrices.

1. LCS Recoveries Criteria

Where LCS spiked with the same analyte at the same concentrations as the MS/MSD? **Yes** or No. If no make note in data review memo.

List the %R of compounds which do not meet the criteria

LCS ID	COMPOUND	% R	QC LIMIT
Recoveries within laboratory control limits except for the cases described in this document.			
GGH5695-BS	Isopropyl Alcohol	125*	76 - 121
	sec-Butyl Alcohol	122*	74 - 118

Note: No action taken, professional judgment. High percent recoveries and no associated positive reported in the QC batch.

* QC limits are laboratory in-house performance criteria, LL = lower limit, UL = upper limit.
* If QC limits are not available, use limits of 70 – 130 %.

Actions:

QUALITY	%R < LL	%R > UL
Positive results	J	J
Nondetects results	R	Accept

All analytes in the associated sample results are qualified for the following criteria.

If 25 % of the LCS recoveries were < LL (or 70 %), qualify all positive results (j) and reject nondetects (R).
If two or more LCS were below 10 %, qualify all positive results as (J) and reject nondetects (R).

2. Frequency Criteria:

Where LCS analyzed at the required frequency and for each matrix? Yes or No.
If no, the data may be affected. Use professional judgment to determine the severity of the effect and qualify data accordingly. Discuss any actions below and list the samples affected.

All criteria were met ___X__
Criteria were not met
and/or see below _____

IX. FIELD/LABORATORY DUPLICATE PRECISION

Sample IDs:__FA42192-8/FA42192-11___ Matrix:___Groundwater_____

Field/laboratory duplicates samples may be taken and analyzed as an indication of overall precision. These analyses measure both field and lab precision; therefore, the results may have more variability than laboratory duplicates which only laboratory performance. It is also expected that Groundwater duplicate results will have a greater variance than water matrices due to difficulties associated with collecting identical field duplicate samples.

The project QAPP should be reviewed for project-specific information.
Suggested criteria: RPD ± 30% for aqueous samples, RPD ± 50 % for solid samples. If both samples and duplicate are <5 SQL, the RPD criteria is doubled.

COMPOUND	SQL	SAMPLE CONC.	DUPLICATE CONC.	RPD	ACTION
Field duplicates analyzed with this data package. RPD within laboratory, generally acceptable and guidance document performance criteria control limits.					

Actions:

Qualify as estimated positive results (J) and nondetects (UJ) for the compound that exceeded the above criteria. For organics, only the sample and duplicate will be qualified.

If an RPD cannot be calculated because one or both of the sample results is not detected, the following actions apply:

If one sample result is not detected and the other is greater than 5x the SQL qualify (J/UJ).

If one sample value is not detected and the other is greater than 5x the SQL and the SQLs for the sample and duplicate are significantly different, use professional judgment to determine if qualification is appropriate.

If one sample value is not detected and the other is less than 5x, use professional judgment to determine if qualification is appropriate.

If both sample and duplicate results are not detected, no action is needed.

All criteria were met __N/A__
Criteria were not met
and/or see below _____

X. INTERNAL STANDARD PERFORMANCE

The assessment of the internal standard (IS) parameter is used to assist the data reviewer in determining the condition of the analytical instrumentation.

List the internal standard area of samples which do not meet the criteria.

* Area of +100% or -50% of the IS area in the associated calibration standard.
* Retention time (RT) within 30 seconds of the IS area in the associated calibration standard.

DATE	SAMPLE ID	IS OUT	IS AREA	ACCEPTABLE RANGE	ACTION

Actions:

1. IS actions should be applied to the compound quantitated with the out-of-control ISs

QUALITY	IS AREA < -25%	IS AREA = -25 % TO – 50%	IS AREA > + 100%
Positive results	J	J	J
Nondetected results	R	UJ	ACCEPT

2. If a IS retention time varies more than 30 seconds, the chromatographic profile for that sample must be examined to determine if any false positive or negative exists. For shifts of a large magnitude, the reviewer may consider partial or total rejection of the data for the sample fraction.

All criteria were met __X__
Criteria were not met
and/or see below _____

XII. SAMPLE QUANTITATION

The sample quantitation evaluation is to verify laboratory quantitation results. In the space below, please show a minimum of one sample calculation:

BS – 03/24/17 Sec-Butyl alcohol RF = 26.42

[ ] = (128619)/(26.42)

= 4,868 ppb OK

All criteria were met __X__
Criteria were not met
and/or see below _____

XII. QUANTITATION LIMITS

A. Dilution performed

SAMPLE ID	DILUTION FACTOR	REASON FOR DILUTION

B. Percent Solids

List samples which have ≤ 50 % solids

Actions:

If the % solids of a Groundwater sample is 10-50%, estimate positive results (J) and nondetects (UJ)

If the % solids of a Groundwater sample is < 10%, estimate positive results (J) and reject nondetects (R)

EXECUTIVE NARRATIVE

SDG No:	**FA42192**	Laboratory:	**Accutest, Orlando**
Analysis:	**MADEP EPH**	Number of Samples:	**11**
Location:	BMSMC, Humacao, PR		

SUMMARY: Eleven (11) samples were analyzed for Extractables TPHC Ranges by method MADEP EPH. Samples were validated following the METHOD FOR THE DETERMINATION OF EXTRACTABLE PETROLEUM HYDROCARBONS (EPH) quality control criteria, Massachusetts Department of Environmental Protection, Revision 1.1 (2004). Also the general validation guidelines promulgated by the USEPA Hazardous Wastes Support Section. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

Results are valid and can be used for decision making purposes.

Critical issues: **None**
Major: **None**
Minor: **None**

Critical findings: **None**
Major findings: **None**
Minor findings: **1.** Ending calibration verification meets method specific requirements except in the cases described in the Data Review Worksheet. C9 – C36 EPH hydrocarbon range were outside method performance criteria. No action taken, professional judgment.

2. C11 – C22 Aromatics (Unadj.) detected in method blank below the reporting limit. Sample results below the reporting limits are qualified as non-detects (U) at the reporting limit.

C11 – C22 Aromatics (Unadj.) detected in field blank below the reporting limit. Sample results below the reporting limits are qualified as non-detects (U) at the reporting limit.

3. No MS/MSD analyzed with this data package. BS/BSD used to assess accuracy. % recovery and RPD within laboratory control limits.

COMMENTS: Results are valid and can be used for decision making purposes.

Reviewers Name: Rafael Infante
Chemist License 1888

Signature: Rafael Infante

Date: April 26, 2017

ORGANIC DATA SAMPLE SUMMARY

Sample ID: FA42192-1
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: AQ - Field Blank Water

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	99.1	ug/L	1.0	J	J	Yes
C9 - C18 Aliphatics	190	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	190	ug/L	1.0	-	U	Yes

Sample ID: FA42192-2
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	200	ug/L	1.0	J	U	Yes
C9 - C18 Aliphatics	200	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	200	ug/L	1.0	-	U	Yes

Sample ID: FA42192-3
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	200	ug/L	1.0	J	U	Yes
C9 - C18 Aliphatics	200	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	200	ug/L	1.0	-	U	Yes

Sample ID: FA42192-6
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	200	ug/L	1.0	J	U	Yes
C9 - C18 Aliphatics	95.3	ug/L	1.0	J	J	Yes
C19 - C36 Aliphatics	50.2	ug/L	1.0	J	J	Yes

Sample ID: FA42192-7
Sample location: BMSMC, Humacao, PR
Sampling date: 3/16/2017
Matrix: Groundwater

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	190	ug/L	1.0	J	U	Yes
C9 - C18 Aliphatics	190	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	190	ug/L	1.0	-	U	Yes

Sample ID: FA42192-8
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	200	ug/L	1.0	J	U	Yes
C9 - C18 Aliphatics	200	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	200	ug/L	1.0	-	U	Yes

Sample ID: FA42192-11
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	200	ug/L	1.0	-	U	Yes
C9 - C18 Aliphatics	200	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	200	ug/L	1.0	-	U	Yes

Sample ID: FA42192-12
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	200	ug/L	1.0	-	U	Yes
C9 - C18 Aliphatics	200	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	200	ug/L	1.0	-	U	Yes

Sample ID: FA42192-13
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	200	ug/L	1.0	J	U	Yes
C9 - C18 Aliphatics	200	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	200	ug/L	1.0	-	U	Yes

Sample ID: FA42192-14
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: Groundwater

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	200	ug/L	1.0	J	U	Yes
C9 - C18 Aliphatics	200	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	200	ug/L	1.0	-	U	Yes

Sample ID: FA42192-15
Sample location: BMSMC, Humacao, PR
Sampling date: 3/17/2017
Matrix: AQ - Field Blank Water

METHOD: MADEP EPH

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
C11 - C22 Aromatics (Unadj.)	200	ug/L	1.0	-	U	Yes
C9 - C18 Aliphatics	200	ug/L	1.0	-	U	Yes
C19 - C36 Aliphatics	200	ug/L	1.0	-	U	Yes

DATA REVIEW WORKSHEETS

Type of validation Full:__X____ Limited:_____

Project Number:_FA42192_____
Date:_____03/16-17/2017_____
Shipping date:___03/17/2017_____
EPA Region:_____2_____

REVIEW OF EXTRACTABLE PETROLEUM HYDROCARBON (EPHs) PACKAGE

The following guidelines for evaluating volatile organics were created to delineate required validation actions. This document will assist the reviewer in using professional judgment to make more informed decision and in better serving the needs of the data users. The sample results were assessed according to the data validation guidance documents in the following order of precedence METHOD FOR THE DETERMINATION OF EXTRACTABLE PETROLEUM HYDROCARBONS (EPH), Massachusetts Department of Environmental Protection, Revision 1.1 (2004). Also the general validation guidelines promulgated by the USEPA Hazardous Wastes Support Section. The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

The hardcopied (laboratory name) _Accutest_Laboratories_____ data package received has been reviewed and the quality control and performance data summarized. The data review for SVOCs included:

Lab. Project/SDG No.: __FA42192_____ Sample matrix: ____Groundater_____
No. of Samples: _____11_____
Field blank No.: _____ FA42192-1;_FA42192-15_____
Equipment blank No.: _-_____
Trip blank No.: _____-_____
Field duplicate No.: ___-_____

__X____ Data Completeness
__X____ Holding Times
__N/A__ GC/MS Tuning
__N/A__ Internal Standard Performance
__X____ Blanks
__X____ Surrogate Recoveries
__X____ Matrix Spike/Matrix Spike Duplicate
___X__ Laboratory Control Spikes
___X__ Field Duplicates
___X__ Calibrations
___X__ Compound Identifications
___X__ Compound Quantitation
___X__ Quantitation Limits

Overall Comments:
_Extractable_Petroleum_Hydrocarbons_by_GC_by_Method_MADEP_EPH,_REV_1.1._____
_____
_____
_____

Definition of Qualifiers:

J- Estimated results
U- Compound not detected
R- Rejected data
UJ- Estimated nondetect

Reviewer:_ Rafael DeFuent _____
Date:_____Apri_25,_2017_____

DATA REVIEW WORKSHEETS

All criteria were met __x___
Criteria were not met and/or see below _____

I. DATA COMPLETNESS
A. Data Package:

MISSING INFORMATION	DATE LAB. CONTACTED	DATE RECEIVED

B. Other Discrepancies:

All criteria were met ___X___
Criteria were not met and/or see below _____

HOLDING TIMES
The objective of this parameter is to ascertain the validity of the results based on the holding time of the sample from time of collection to the time of extraction, and subsequently from the time of extraction to the time of analysis.

Complete table for all samples and note the analysis and/or preservation not within criteria

SAMPLE ID	DATE SAMPLED	DATE EXTRACTED	DATE ANALYZED	ACTION
Samples extracted and analyzed within method recommended holding time				

Criteria

Preservation:

Aqueous samples must be acidified to a pH of 2.0 or less at the time of collection.
Soil samples must be cooled at 4 ± 2 °C immediately after collection.

Holding times:

Samples must be extracted within 14 days of collection, and analyzed within 40 days of extraction.

Cooler temperature (Criteria: 4 ± 2 °C):___1.6_-_5.8_°C_____

Actions: Qualify positive results/nondetects as follows:

If holding times are exceeded, estimate positive results (J) and nondetects (UJ).
If holding times are grossly exceeded, use professional judgment to qualify data. The data reviewer may choose to estimate positive results (J) and rejects nondetects (R).
If samples were not at the proper temperature (> 10°C) or improperly preserved, use professional judgment to qualify the results.

Note:

All criteria were met ___X___
Criteria were not met and/or see below _____

CALIBRATIONS VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:_____03/15/17;_04/10/17_____

Dates of initial calibration verification:_____03/15/16;_04/10/17_____

Instrument ID numbers:_____FID_7_____

Matrix/Level:_____AQUEOUS/MEDIUM_____

DATE	LAB FILE ID#	ANALYTE	CRITERIA OUT RFs, %RSD, %D, r	SAMPLES AFFECTED
Initial and initial calibration verification meet method specific requirements				

Criteria- ICAL

- Five point calibration curve.
- The percent relative standard deviation (%RSD) of the calibration factor must be equal to or less than 25% over the working range for the analyte of interest. When this condition is met, linearity through the origin may be assumed, and the average calibration factor is used in lieu of a calibration curve.
- A collective calibration factor must also be established for each hydrocarbon range of interest. Calculate the collective CFs for C9-C18 Aliphatic Hydrocarbons, C19-C36 Aliphatic Hydrocarbons, and C11-C22 Aromatic Hydrocarbons using the FID chromatogram. Tabulate the summation of the peak areas of all components in that fraction against the total concentration injected. The %RSD of the calibration factor must be equal to or less than 25% over the working range for the hydrocarbon range of interest.
    - The area for the surrogates must be subtracted from the area summation of the range in which they elute.
    - The areas associated with naphthalene and 2-methylnaphthalene in the aliphatic range standard must be subtracted from the uncorrected collective C9-C18 Aliphatic Hydrocarbon range area prior to calculating the CF.

Criteria- CCAL

- At a minimum, the working calibration factor must be verified on each working day, after every 20 samples or every 24 hours (whichever is more frequent), and at the end of the analytical sequence by the injection of a mid-level continuing calibration standard to verify instrument performance and linearity.
- If the percent difference (%D) for any analyte varies from the predicted response by more than ±25%, a new five-point calibration must be performed for that analyte. Greater percent differences are permissible for n-nonane. If the %D for n-nonane is greater than 30, note the nonconformance in the case narrative. It should be noted that the %Ds are calculated when CFs are used for the initial calibration and percent drifts are calculated when calibration curves using linear regression are used for the initial calibration.

Actions:

If %RSD > 25% for target compounds or a correlation coefficient < 0.99, estimate positive results (J) and use professional judgment to qualify nondetects.
If % D > 25% (> 30 for nonane), estimate positive results (J) and nondetects (UJ).

CALIBRATIONS VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:_____04/10/17_____

Dates of continuing calibration verification:___04/12/17;_04/13/17;_04/15/17;_04/17/17_

Dates of final calibration verification:_____04/14/17;_04/15/17;_04/17/17_____

Instrument ID numbers:_____FID_7_____

Matrix/Level:_____AQUEOUS/MEDIUM_____

DATE	LAB FILE ID#	ANALYTE	CRITERIA OUT RFs, %RSD, %D, r	SAMPLES AFFECTED
Continuing calibration meets method specific requirements except in the cases described in this document. Ending calibration verification included in data package.				
04/14/17	ecc911-4/ ecc898-4	C9 – C36 Aliphatics	-25.6	FA42192-13; -14
04/17/17	ecc911-4/ ecc898-4	C9 – C36 Aliphatics	-27.3	FA42192-12

Note: No action taken, professional judgment.

A separate worksheet should be filled for each initial curve

All criteria were met _____
Criteria were not met and/or see below ___X___

V A. BLANK ANALYSIS RESULTS (Sections 1 & 2)

The assessment of the blank analysis results is to determine the existence and magnitude of contamination problems. The criteria for evaluation of blanks apply only to blanks associated with the samples, including trip, equipment, and laboratory blanks. If problems with any blanks exist, all data associated with the case must be carefully evaluated to determine whether or not there is an inherent variability in the data for the case, or if the problem is an isolated occurrence not affecting other data. A Laboratory Method Blank must be run after samples suspected of being highly contaminated to determine if sample carryover has occurred.

List the contamination in the blanks below. High and low levels blanks must be treated separately.

Laboratory blanks

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
METHOD BLANKS MEET THE METHOD SPECIFIC CRITERIA EXCEPT FOR THE CASES DESCRIBED IN THIS DOCUMENT.				
04/13/17	OP64317-MB	Aq./medium	C11-C22 Aromatics (Unadj.)	87.2 ug/l

Note: Blank concentration below reporting limits. Sample results below the reporting limits are qualified as non-detects (U) at the reporting limit.

Field/Trip/Equipment

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
NO TARGET ANALYTES DETECTED IN THE FIELD BLANKS ANALYZED WITH THIS DATA PACKAGE EXCEPT IN THE CASES DESCRIBED IN THIS DOCUMENT. NO EQUIPMENT BLANK ANALYZED ASSOCIATED WITH THIS DATA PACKAGE.				
04/12/17	FA42192-1	Aq./medium	C11-C22 Aromatics (Unadj.)	99.1 ug/l

Note: Blank concentration below reporting limits. Sample results below the reporting limits are qualified as non-detects (U) at the reporting limit.

All criteria were met _____
Criteria were not met and/or see below __X___

V B. BLANK ANALYSIS RESULTS (Section 3)

Blank Actions

The ALs for samples which have been diluted should be corrected for the sample dilution factor and/or % moisture, where applicable. Peaks must not be detected above the Reporting Limit within the retention time window of any analyte of interest. The hydrocarbon ranges must not be detected at a concentration greater than 10% of the most stringent MCP cleanup standard. Specific actions area as follows:

If the concentration is < sample quantitation limit (SQL) and < AL, report the compound as not detected (U) at the SQL.
If the concentration is ≥ SQL but < AL, report the compound as not detected (U) at the reported concentration.
If the concentration is > AL, report the concentration unqualified.

All criteria were met __X___
Criteria were not met and/or see below _____

SURROGATE SPIKE RECOVERIES

Laboratory performance of individual samples is established by evaluation of surrogate spike recoveries. All samples are spiked with surrogate compounds prior to sample analysis. The accuracy of the analysis is measured by the surrogate percent recovery. Since the effects of the sample matrix are frequently outside the control of the laboratory and may present relatively unique problems, the validation of data is frequently subjective and demands analytical experience and professional judgment.
List the percent recoveries (%Rs) which do not meet the criteria for surrogate recovery.
Matrix: solid/aqueous

SAMPLE ID	**SURROGATE COMPOUND**				**ACTION**
	S1	S2	S3	S4	

_SURROGATE_STANDARDS_RECOVERIES_WITHIN_LABORATORY_CONTROL__
_LIMITS._____
_____
_____

Note:

S1 = o-Terphenyl 40-140% S2 = 2-Fluorobiphenyl 40-140%
S3 = 1-Chlorooctadecane 40-140% S4 = 2-Bromonaphthalene 40-140%

QC Limits (%)* (Aqueous)
_LL_to_UL_ _40_to_140_ _40_to_140_ _40_to_140_ _40_to_140_
QC Limits* (Solid)
_LL_to_UL_ ___to___ ____to_____ ___to_____ ___to_____

It is recommended that surrogate standard recoveries be monitored and documented on a continuing basis. At a minimum, when surrogate recovery from a sample, blank, or QC sample is less than 40% or more than 140%, check calculations to locate possible errors, check the fortifying standard solution for degradation, and check changes in instrument performance.

If the cause cannot be determined, reanalyze the sample unless one of the following exceptions applies:

(1) Obvious interference is present on the chromatogram (e.g., unresolved complex mixture);
(2) The surrogate exhibits high recovery and associated target analytes or hydrocarbon ranges are not detected in sample.

If a sample with a surrogate recovery outside of the acceptable range is not reanalyzed based on any of these aforementioned exceptions, this information must be noted on the data report form and discussed in the Executive Report. Analysis of the sample on dilution may diminish matrix-related surrogate recovery problems. This approach can be used as long as the reporting limits to evaluate applicable MCP standards can still be achieved with the dilution. If not, reanalysis without dilution must be performed.

All criteria were met _____
Criteria were not met and/or see below __N/A__

VII. A MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

This data is generated to determine long term precision and accuracy in the analytical method for various matrices. This data alone cannot be used to evaluate the precision and accuracy of individual samples.

At the request of the data user, and in consideration of sample matrices and data quality objectives, matrix spikes and matrix duplicates may be analyzed with every batch of 20 samples or less per matrix.

- **Matrix duplicate** - Matrix duplicates are prepared by analyzing one sample in duplicate. The purpose of the matrix duplicates is to determine the homogeneity of the sample matrix as well as analytical precision. The RPD of detected results in the matrix duplicate samples must not exceed 50 when the results are greater than 5x the reporting limit.
- The desired spiking level is 50% of the highest calibration standard. However, the total concentration in the MS (including the MS and native concentration in the unspiked sample) should not exceed 75% of the highest calibration standard in order for a proper evaluation to be performed. The purpose of the matrix spike is to determine whether the sample matrix contributes bias to the analytical results. The corrected concentrations of each analyte within the matrix spiking solution must be within 40 - 140% of the true value. Lower recoveries of n-nonane are permissible but must be noted in the narrative if <30%.

MS/MSD Recoveries and Precision Criteria

Sample ID:_____-_____ Matrix/Level:_____-_____

List the %Rs, RPD of the compounds which do not meet the QC criteria.

MS OR MSD	COMPOUND	% R	RPD	QC LIMITS	ACTION

Note: No MS/MSD analyzed with this data package. BS/BSD used to assess accuracy. % recovery and RPD within laboratory control limits.

All criteria were met _____
Criteria were not met and/or see below __N/A__

No action is taken on MS/MSD results alone to qualify the entire case. However, used informed professional judgment, the data reviewer may use the MS/MSD results in conjunction with other QC criteria and determine the need for some qualification of the data. In those instances where it can be determined that the results of the MS/MSD affect only the sample spiked, the qualification should be limited to this sample alone. However, it may be determined through the MS/MSD results that the laboratory is having a systematic problem in the analysis of one or more analytes, which affects the associated samples.

2. MS/MSD – Unspiked Compounds

List the concentrations of the unspiked compounds and determine the % RSDs of these compounds in the unspiked sample, matrix spike, and matrix spike duplicate.

COMPOUND	CONCENTRATION SAMPLE	MS	MSD	%RPD	ACTION

Criteria: None specified, use %RSD $\leq$ 50 as professional judgment.

Actions:

If the % RSD > 50, qualify the results in the spiked sample as estimate (J).
If the % RSD is not calculable (NC) due to nondetect value in the sample, MS, and/or MSD, use professional judgment to qualify sample data.

A separate worksheet should be used for each MS/MSD pair.

All criteria were met __X___
Criteria were not met and/or see below _____

VIII. LABORATORY CONTROL SAMPLE (LCS/LCSD) ANALYSIS

This data is generated to determine accuracy of the analytical method for various matrices.

1. LCS Recoveries Criteria

List the %R of compounds which do not meet the criteria

LCS ID	COMPOUND	% R	QC LIMIT	ACTION

__LCS/LCSD_RECOVERY_WITHIN_LABORATORY_CONTROL_LIMTS._____

Criteria:
* Refer to QAPP for specific criteria.
* The spike recovery must be between 40% and 140%. Lower recoveries of n-nonane are permissible. If the recovery of n-nonane is <30%, note the nonconformance in the executive narrative. RPD between LCS/LCSD must be < 25%.

Actions:
Actions on LCS recovery should be based on both the number of compounds that are outside the %R and RPD criteria and the magnitude of the excedance of the criteria.

If the %R of the analyte is > UL, qualify all positive results (j) for the affected analyte in the associated samples and accept nondetects.
If the %R of the analyte is < LL, qualify all positive results (j) and reject (R) nondetects for the affected analyte in the associated samples.
If more than half the compounds in the LCS are not within the required recovery criteria, qualify all positive results as (J) and reject nondetects (R) for all target analyte(s) in the associated samples.

2. Frequency Criteria:

Where LCS analyzed at the required frequency and for each matrix (1 per 20 samples per matrix)? **Yes** or No.
If no, the data may be affected. Use professional judgment to determine the severity of the effect and qualify data accordingly. Discuss any actions below and list the samples affected. Discuss the actions below:

All criteria were met __X___
Criteria were not met and/or see below ___

IX. FIELD/LABORATORY DUPLICATE PRECISION

Sample IDs: __FA42192-8/-11_(field)_____ Matrix:____Groundwater___
Sample IDs: __FA42192-7/-7DUP_(lab)_____ Matrix:____Groundwater___

Field/laboratory duplicates samples may be taken and analyzed as an indication of overall precision. These analyses measure both field and lab precision; therefore, the results may have more variability than laboratory duplicates which measures only laboratory performance. It is also expected that soil duplicate results will have a greater variance than water matrices due to difficulties associated with collecting identical field duplicate samples.

COMPOUND	SQL	SAMPLE CONC.	DUPLICATE CONC.	RPD	ACTION
Field/laboratory duplicate analyzed with this data package. RPD within laboratory and generally acceptable control limits for analytes detected above 5 x SQL.					

Criteria:

The project QAPP should be reviewed for project-specific information.
RPD $\pm$ 30% for aqueous samples, RPD $\pm$ 50 % for solid samples if results are $\geq$ SQL.
If both samples and duplicate are <5 SQL, the RPD criteria is doubled.

SQL = soil quantitation limit

Actions:

If both the sample and the duplicate results are nondetects (ND), the RPD is not calculable (NC). No action is needed.

Qualify as estimated positive results (J) and nondetects (UJ) for the compound that exceeded the above criteria.

If one sample result is not detected and the other is $\geq$ 5x the SQL qualify (J/UJ).

Note: If SQLs for the sample and duplicate are significantly different, use professional judgment to determine if qualification is appropriate.

If one sample value is not detected and the other is < 5x the SQL, use professional judgment to determine if qualification is appropriate.

All criteria were met __X___
Criteria were not met and/or see below _____

XI. COMPOUND IDENTIFICATION

The compound identification evaluation is to verify that the laboratory correctly identified target analytes as well as tentatively identified compounds (TICs).

1. Verify that the target analytes were within the retention time windows.

   - Retention time windows must be re-established for each Target EPH Analyte each time a new GC column is installed, and must be verified and/or adjusted on a daily basis.
   - The n-nonane (n-C9) peak must be adequately resolved from the solvent front of the chromatographic run.
   - All surrogates must be adequately resolved from the Aliphatic Hydrocarbon and Aromatic Hydrocarbon standards.
   - For the purposes of this method, adequate resolution is assumed to be achieved if the height of the valley between two peaks is less than 25% of the average height of the two peaks.
   - The n-pentane (C5) and MtBE peaks must be adequately resolved from any solvent front that may be present on the FID and PID chromatograms, respectively.

1a. Aliphatic hydrocarbons range:

   - Determine the total area count for all peaks eluting 0.1 minutes before the retention time (Rt) for n-C9 and 0.01 minutes before the Rt for n-C19.
   - Determine the total area count for all peaks eluting 0.01 minutes before the Rt for n-C19 and 0.1 minutes after the Rt for n-C36.

Are the aliphatic hydrocarbons range properly determined? **Yes**? or No?

Comments:

1b. Aromatic hydrocarbons range:

   - Determine the total area count for all peaks eluting 0.1 minutes before the retention time (Rt) for naphthalene and 0.1 minutes after the Rt for benzo(g,h,i)perylene.
   - Determine the peak area count for the sample surrogate (OTP) and fractionation surrogate(s). Subtract these values from the collective area count value.

Are the aliphatic hydrocarbons range properly determined? **Yes**? or No?

Comments:

All criteria were met __X___
Criteria were not met and/or see below _____

2. If target analytes and/or TICs were not correctly identified, request that the laboratory resubmit the corrected data.

3. Breakthrough determination - Each sample (field and QC sample) must be evaluated for potential breakthrough on a sample specific basis by evaluating the % recovery of the fractionation surrogate (2-bromonaphthalene) and on a batch basis by quantifying naphthalene and 2-methylnaphthalene in both the aliphatic and aromatic fractions of the LCS and LCSD. **If either the concentration of naphthalene or 2-methylnaphthalene in the aliphatic fraction exceeds 5% of the total concentration for naphthalene or 2-methylnaphthalene in the LCS or LCSD, fractionation must be repeated on all archived batch extracts.**

NOTE: **The total concentration of naphthalene or 2-methylnaphthalene in the LCS/LCSD pair includes the summation of the concentration detected in the aliphatic fraction and the concentration detected in the aromatic fraction.**

Comments:__Concentration_in_the_aliphatic_fraction_<_5%_of_the_total_____
_concentration_for_naphthalene_and_2-methylnaphthalene_____
_____
_____
_____

4. **Fractionation Check Standard** – A fractionation check solution is prepared containing 14 alkanes and 17 PAHs at a nominal concentration of 200 ng/µl of each constituent. The Fractionation Check Solution must be used to evaluate the fractionation efficiency of each new lot of silica gel/cartridges, and establish the optimum hexane volume required to efficiently elute aliphatic hydrocarbons while not allowing significant aromatic hydrocarbon breakthrough. For each analyte contained in the fractionation check solution, excluding n-nonane, the Percent Recovery must be between 40 and 140%. A 30% Recovery is acceptable for n-nonane.

Is a fractionation check standard analyzed? Yes? or No?

Comments: Not applicable.

All criteria were met __X___
Criteria were not met and/or see below _____

XII. QUANTITATION LIMITS AND SAMPLE RESULTS

The sample quantitation evaluation is to verify laboratory quantitation results.

In order to demonstrate the absence of aliphatic mass discrimination, the response ratio of C28 to C20 must be at least 0.85. If <0.85, this nonconformance must be noted in the laboratory case narrative.

The chromatograms of Continuing Calibration Standards for aromatics must be reviewed to ensure that there are no obvious signs of mass discrimination.

Is aliphatic mass discrimination observed in the sample? Yes? or **No**?

Is aromatic mass discrimination observed in the sample? Yes? or **No**?

1. In the space below, please show a minimum of one sample calculation:

   FA42192-6 (C9 – C18, Aliphatics) RF = 1.074×10^6

   [ ] = $51147395/1.074 \times 10^6$

   = 47.62 ppm

2. If requested, verify that the results were above the laboratory method detection limit (MDLs).

3. If dilutions performed, were the SQLs elevated accordingly by the laboratory? List the affected samples and dilution factor in the table below.

SAMPLE ID	DILUTION FACTOR	REASON FOR DILUTION

If dilution was not performed, estimate results (J) for the affected compounds. List the affected samples/compounds:

_____

_____

EXECUTIVE NARRATIVE

SDG No:	**FA42192**	Laboratory:	**Accutest, Orlando**
Analysis:	**SW846-8081B**	Number of Samples:	**13**

Location: BMSMC, Humacao, PR

SUMMARY: Thirteen (13) samples were analyzed for the TCL pesticides list following method SW846-8081B. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence *Hazardous Waste Support Section SOP No. HW-36A, Revision 0, June, 2015. SOM02.2. Pesticide Data Validation.* The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

Results are valid and can be used for decision making purposes.

Critical issues: **None**
Major: **None**
Minor: **None**

Critical findings: **None**
Major findings: **None**
Minor findings:

1. Initial and initial calibration verification within the guidance document performance criteria. Continuing calibration % differences meet the performance criteria in at least one of the two columns except for the cases described in the Data Review Worksheet. Results for Endrin ketone qualified as estimated (J or UJ) in affected samples.

Several analytes missed the % difference continuing calibration criteria in cc1931-40 analyzed on 03/28/17 (9:49:54 PM). No samples from this batch associated with this continuing calibration check.

Final calibration verification included in data package.

2. Surrogate recoveries were within laboratory control limits except in the cases described in the Data Review Worksheet. No action taken, professional judgment. Surrogates recovered high and analytes not detected in affected samples.

3. MS/MSD % recoveries and RPD within laboratory control limits except for the cases described in the Data Review Worksheet. No qualification made professional judgment. Analytes recovered high and not detected in sample batch above reporting limit except for alpha-Chlordane.

MS/MSD % recovery results apply to unspiked sample. Positive alpha-Chlordane results were from samples FA42192-7 and FA42192-12. No qualification made.

4. LCS % recovery and RPD within laboratory control limits except in the cases described in the Data Review Worksheet. LCS recovery exceeds the upper acceptance limit; non-detected compounds are not qualified.

COMMENTS: Results are valid and can be used for decision making purposes.

Reviewers Name: Rafael Infante
Chemist License 1888

Signature: Rafael Infante

Date: April 26, 2017

SAMPLE ORGANIC DATA SAMPLE SUMMARY

Sample ID: FA42192-1
Sample location: BMSMC, Humacao, PR
Sampling date: 16-Mar-17
Matrix: AQ - Fielb Blank Water

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.040	ug/l	1	-	U	Yes
gamma-Chlordane	0.040	ug/l	1	-	U	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-2
Sample location: BMSMC, Humacao, PR
Sampling date: 16-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.040	ug/l	1	-	U	Yes
gamma-Chlordane	0.013	ug/l	1	J	J	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-3
Sample location: BMSMC, Humacao, PR
Sampling date: 16-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.040	ug/l	1	-	U	Yes
gamma-Chlordane	0.017	ug/l	1	J	J	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-6
Sample location: BMSMC, Humacao, PR
Sampling date: 16-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.024	ug/l	1	J	J	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.040	ug/l	1	-	U	Yes
gamma-Chlordane	0.040	ug/l	1	-	U	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	U	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-7
Sample location: BMSMC, Humacao, PR
Sampling date: 16-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.041	ug/l	1	-	-	Yes
gamma-Chlordane	0.040	ug/l	1	-	U	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-8
Sample location: BMSMC, Humacao, PR
Sampling date: 17-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.040	ug/l	1	-	U	Yes
gamma-Chlordane	0.040	ug/l	1	-	U	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-11
Sample location: BMSMC, Humacao, PR
Sampling date: 17-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.040	ug/l	1	-	U	Yes
gamma-Chlordane	0.040	ug/l	1	-	U	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-12
Sample location: BMSMC, Humacao, PR
Sampling date: 17-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.046	ug/l	1	-	-	Yes
gamma-Chlordane	0.040	ug/l	1	-	U	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-13
Sample location: BMSMC, Humacao, PR
Sampling date: 17-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.025	ug/l	1	J	J	Yes
gamma-Chlordane	0.040	ug/l	1	-	U	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-14
Sample location: BMSMC, Humacao, PR
Sampling date: 17-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.040	ug/l	1	-	U	Yes
gamma-Chlordane	0.040	ug/l	1	-	U	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-15
Sample location: BMSMC, Humacao, PR
Sampling date: 17-Mar-17
Matrix: AQ - Field Blank Water

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	0.040	ug/l	1	-	U	Yes
alpha-BHC	0.040	ug/l	1	-	U	Yes
beta-BHC	0.040	ug/l	1	-	U	Yes
delta-BHC	0.040	ug/l	1	-	U	Yes
gamma-BHC (Lindane)	0.040	ug/l	1	-	U	Yes
alpha-Chlordane	0.040	ug/l	1	-	U	Yes
gamma-Chlordane	0.040	ug/l	1	-	U	Yes
Dieldrin	0.040	ug/l	1	-	U	Yes
4,4'-DDD	0.080	ug/l	1	-	U	Yes
4,4'-DDE	0.080	ug/l	1	-	U	Yes
4,4'-DDT	0.080	ug/l	1	-	U	Yes
Endrin	0.080	ug/l	1	-	U	Yes
Endosulfan sulfate	0.080	ug/l	1	-	U	Yes
Endrin aldehyde	0.080	ug/l	1	-	U	Yes
Endrin ketone	0.080	ug/l	1	-	UJ	Yes
Endosulfan-I	0.040	ug/l	1	-	U	Yes
Endosulfan-II	0.040	ug/l	1	-	U	Yes
Heptachlor	0.040	ug/l	1	-	U	Yes
Heptachlor epoxide	0.040	ug/l	1	-	U	Yes
Methoxychlor	0.080	ug/l	1	-	U	Yes
Toxaphene	2.0	ug/l	1	-	U	Yes

Sample ID: FA42192-3 MS
Sample location: BMSMC, Humacao, PR
Sampling date: 16-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	1.2	ug/l	1	-	-	Yes
alpha-BHC	1.2	ug/l	1	-	-	Yes
beta-BHC	1.2	ug/l	1	-	-	Yes
delta-BHC	1.3	ug/l	1	-	-	Yes
gamma-BHC (Lindane)	1.2	ug/l	1	-	-	Yes
alpha-Chlordane	1.4	ug/l	1	-	-	Yes
gamma-Chlordane	1.7	ug/l	1	-	-	Yes
Dieldrin	1.3	ug/l	1	-	-	Yes
4,4'-DDD	1.4	ug/l	1	-	-	Yes
4,4'-DDE	1.3	ug/l	1	-	-	Yes
4,4'-DDT	1.4	ug/l	1	-	-	Yes
Endrin	1.4	ug/l	1	-	-	Yes
Endosulfan sulfate	1.4	ug/l	1	-	-	Yes
Endrin aldehyde	1.3	ug/l	1	-	-	Yes
Endrin ketone	1.5	ug/l	1	-	-	Yes
Endosulfan-I	1.2	ug/l	1	-	-	Yes
Endosulfan-II	1.3	ug/l	1	-	-	Yes
Heptachlor	1.2	ug/l	1	-	-	Yes
Heptachlor epoxide	1.2	ug/l	1	-	-	Yes
Methoxychlor	1.4	ug/l	1	-	-	Yes

Sample ID: FA42192-3 MSD
Sample location: BMSMC, Humacao, PR
Sampling date: 16-Mar-17
Matrix: Groundwater

METHOD: 8081B

Analyte Name	Result	Units	Dilution Factor	Lab Flag	Validation	Reportable
Aldrin	1.3	ug/l	1	-	-	Yes
alpha-BHC	1.3	ug/l	1	-	-	Yes
beta-BHC	1.3	ug/l	1	-	-	Yes
delta-BHC	1.4	ug/l	1	-	-	Yes
gamma-BHC (Lindane)	1.3	ug/l	1	-	-	Yes
alpha-Chlordane	1.5	ug/l	1	-	-	Yes
gamma-Chlordane	2.0	ug/l	1	-	-	Yes
Dieldrin	1.4	ug/l	1	-	-	Yes
4,4'-DDD	1.5	ug/l	1	-	-	Yes
4,4'-DDE	1.4	ug/l	1	-	-	Yes
4,4'-DDT	1.5	ug/l	1	-	-	Yes
Endrin	1.5	ug/l	1	-	-	Yes
Endosulfan sulfate	1.5	ug/l	1	-	-	Yes
Endrin aldehyde	1.3	ug/l	1	-	-	Yes
Endrin ketone	1.5	ug/l	1	-	-	Yes
Endosulfan-I	1.2	ug/l	1	-	-	Yes
Endosulfan-II	1.4	ug/l	1	-	-	Yes
Heptachlor	1.3	ug/l	1	-	-	Yes
Heptachlor epoxide	1.3	ug/l	1	-	-	Yes
Methoxychlor	1.5	ug/l	1	-	-	Yes

DATA REVIEW WORKSHEETS

Project/CasNumber:_____FA42192_____
Sampling Date:_____03/16-14/2017_____
Shipping Date:_____03/17/2017_____
EPA Region No.:_____2_____

REVIEW OF PESTICIDE ORGANIC PACKAGE

The following guidelines for evaluating volatile organics were created to delineate required validation actions. This document will assist the reviewer in using professional judgment to make more informed decision and in better serving the needs of the data users. The sample results were assessed according to USEPA data validation guidance documents in the following order of precedence *Hazardous Waste Support Section SOP No. HW-36A, Revision 0, June, 2015. SOM02.2. Pesticide Data Validation.* The QC criteria and data validation actions listed on the data review worksheets are from the primary guidance document, unless otherwise noted.

The hardcopied (laboratory name) _Accutest_____ data package received has been reviewed and the quality control and performance data summarized. The data review for VOCs included:

Lab. Project/SDG No.: __FA42192_____ Sample matrix: __Groundwater_
No. of Samples: _____13_____

Trip blank No.: _____-_____
Field blank No.: _____ FA42192-1;_ FA42192-15_____
Equipment blank No.:_____-_____
Field duplicate No.:_____-_____
Field spikes No.:_____FA42192-3MS/-3MSD_____
QC audit samples:_____-_____

___X___ Data Completeness
___X___ Holding Times
__N/A__ GC/MS Tuning
___X___ Internal Standard Performance
___X___ Blanks
___X___ Surrogate Recoveries
___X___ Matrix Spike/Matrix Spike Duplicate
__X____ Laboratory Control Spikes
__X____ Field Duplicates
__X____ Calibrations
__X____ Compound Identifications
__X____ Compound Quantitation
__X____ Quantitation Limits

Overall Comments:__Dieldrin_by_SW846-8081B_____

Definition of Qualifiers:

J-	Estimated results	U-	Compound not detected
R-	Rejected data	UJ-	Estimated nondetect

Reviewer:_____
Date:__April_25,_2017_____

DATA REVIEW WORKSHEETS

DATA COMPLETENESS

MISSING INFORMATION	DATE LAB. CONTACTED	DATE RECEIVED

All criteria were met __X__
Criteria were not met
and/or see below _____

HOLDING TIMES

The objective of this parameter is to ascertain the validity of the results based on the holding time of the sample from time of collection to the time of analysis.

Complete table for all samples and note the analysis and/or preservation not within criteria

SAMPLE ID	DATE SAMPLED	DATE EXTRACTED/ANALYZED	ACTION
Samples properly preserved. All samples extracted and analyzed within the required criteria.			

Note:

Criteria

Aqueous samples - seven (7) days from sample collection for extraction; 40 days from sample collection for analysis.
Non-aqueous samples – fourteen (14) days from sample collection for extraction; 40 days from sample collection for analysis.

Cooler temperature (Criteria: 4 ± 2 °C): 1.6 – 5.8 °C - OK

Actions

Qualify aqueous sample results using preservation and technical holding time information as follows:

a. If there is no evidence that the samples were properly preserved (T = 4°C ± 2°C), and the samples were extracted or analyzed within the technical holding times, qualify detects as estimated (J) and non-detects as estimated (UJ).

b. If there is no evidence that the samples were properly preserved (T = 4°C ± 2°C), and the samples were extracted or analyzed outside the technical holding times, qualify detects as estimated (J) and non-detects as estimated (UJ).

c. If the samples were properly preserved, and were extracted and analyzed within the technical holding times, no qualification of the data is necessary.

d. If the samples were properly preserved, and were extracted or analyzed outside the technical holding times, qualify detects as estimated (J) and non-detects as estimated (UJ). Note in the Data Review Narrative that holding times were exceeded and the effect of exceeding the holding time on the resulting data.

e. Use professional judgment to qualify samples whose temperature upon receipt at the laboratory is either below 2 degrees centigrade or above 6 degrees centigrade.
f. If technical holding times are grossly exceeded, use professional judgment to qualify the data.

Qualify non-aqueous sample results using preservation and technical holding time information as follows:
a. If there is no evidence that the samples were properly preserved (T = 4°C ± 2°C), and the samples were extracted or analyzed within the technical holding time, qualify detects as estimated (J) and non-detects as estimated (UJ).
b. If there is no evidence that the samples were properly preserved (T = 4°C ± 2°C), and the samples were extracted or analyzed outside the technical holding time, qualify detects as estimated (J) and non-detects as estimated (UJ).
c. If the samples were properly preserved, and were extracted and analyzed within the technical holding time, no qualification of the data is necessary.
d. If the samples were properly preserved, and were extracted or analyzed outside the technical holding time, qualify detects as estimated (J) and non-detects as estimated (UJ). Note in the Data Review Narrative that holding times were exceeded and the effect of exceeding the holding time on the resulting data.
e. Use professional judgment to qualify samples whose temperature upon receipt at the laboratory is either below 2 degrees centigrade or above 6 degrees centigrade.
f. If technical holding times are grossly exceeded, use professional judgment to qualify the data.

All criteria were met ___X___
Criteria were not met see below _____

GAS CHROMATOGRAPH WITH ELECTRON CAPTURE DETECTOR (GC/ECD) INSTRUMENT PERFORMANCE CHECK (SECTIONS 1 TO 5)

1. Resolution Check Mixture

Criteria

Is the resolution between two adjacent peaks in the Resolution Check Mixture C greater than or equal to 80.0% for all analytes for the primary column and greater than or equal to 50.0% for the confirmation column? **Yes**? or No?

Is the resolution between two adjacent peaks in the Resolution Check Mixture (A and B) greater than or equal to 60.0%? Yes? or No?

Note: If resolution criteria are not met, the quantitative results may not be accurate due to inadequate resolution. Qualitative identifications may also be questionable if coelution exists.

Action

a. Qualify detects for target compounds that were not adequately resolved as tentatively identified (NJ).
b. Qualify non-detected compounds as unusable (R).

2. Performance Evaluation Mixture (PEM) Resolution Criteria

Criteria

Is PEM analysis performed at the required frequency (at the end of each pesticide initial calibration sequence and every 12 hours)? **Yes**? or No?

Action

a. If PEM is not performed at the required frequency, qualify all associated sample and blank results as unusable (R).

Criteria

Is PEM % Resolution < 90%? Yes? or **No?**

Action

a. a. Qualify detects for target compounds that were not adequately resolved as tentatively identified (NJ).
b. Qualify non-detected compounds as unusable (R).

All criteria were met ___X___
Criteria were not met see below _____

3. PEM 4,4'-DDT Breakdown

Criteria

Is the PEM 4,4'-DDT % Breakdown >20.0% and 4,4'-DDT is detected? Yes? or **No**?

Action

a. Qualify detects for 4,4'-DDT; detects for 4,4'-DDD; and detects for 4,4'-DDE as estimated (J)

Criteria

Is the PEM 4,4'-DDT % Breakdown >20.0% and 4,4'-DDT is not detected Yes? or **No**?

Action

a. Qualify non-detects for 4,4'- DDT as unusable (R)
b. Qualify detects for 4,4'-DDD as tentatively identified (NJ)
c. Qualify detects for 4,4'-DDE as tentatively identified (NJ)

4. PEM Endrin Breakdown

Criteria

Is the PEM Endrin % Breakdown >20.0% and Endrin is detected? Yes? or **No**?

Action

a. Qualify detects for Endrin; detects for Endrin aldehyde; and detects for Endrin ketone as estimated (J)

Criteria

Is the PEM Endrin % Breakdown >20.0% and Endrin is not detected Yes? or **No**?

Action

a. Qualify non-detects for Endrin as unusable (R)
b. Qualify detects for Endrin aldehyde as tentatively identified (NJ)
c. Qualify detects for Endrin ketone as tentatively identified (NJ)

All criteria were met ___X___
Criteria were not met see below _____

5. Mid-point Individual Standard Mixture Resolution -

Criteria

Is the resolution between two adjacent peaks in the Resolution Check Mixture C greater than or equal to 80.0% for all analytes for the primary column and greater than or equal to 50.0% for the confirmation column? **Yes**? or No?

Is the resolution between two adjacent peaks in the Resolution Check Mixture (A and B) greater than or equal to 90.0%? **Yes**? or No?

Note: If resolution criteria are not met, the quantitative results may not be accurate due to inadequate resolution. Qualitative identifications may also be questionable if coelution exists.

Action

a. Qualify detects for target compounds that were not adequately resolved as tentatively identified (NJ).
b. Qualify non-detected compounds as unusable (R).

Criteria

Is mid-point individual standard mixture analysis performed at the required frequency (every 12 hours)? **Yes**? or No?

Action

a. If the mid-point individual standard mixture analysis is not performed at the required frequency, qualify all associated sample and blank results as unusable (R).

All criteria were met ___X___
Criteria were not met
and/or see below _____

CALIBRATION VERIFICATION

Compliance requirements for satisfactory instrument calibration are established to ensure that the instrument is capable of producing and maintaining acceptable quantitative data.

Date of initial calibration:_____04/03/17_____ _____03/24/17_____
Dates of initial calibration verification:_____04/03/17_____ _____03/24/17_____
Dates of continuing calibration:_____04/03/17_____ _____03/28/17_____
Dates of final calibration_____-_____ ___03/24/17;_03/29/17__
Instrument ID numbers:_____ECD_5_____ ____ECD_6_____
Matrix/Level:_____Aqueous/low_____ ___Aqueous/low_____

DATE	LAB FILE ID#	CRITERIA OUT RFs, %RSD, **%D**, r	COMPOUND	SAMPLES AFFECTED
03/28/17 (4:43 PM)	cc1931-40	-20.8/-26.5	Endrin ketone	FA42192-1; -2; -3 ✓
03/28/17 (7:39 PM)	cc1931-40	-21.7/-22.9	Endrin ketone	FA42192-7; -8; -11; -12; -13; -14; -15 ✓

Note: Initial and initial calibration verification within the guidance document performance criteria. Continuing calibration % differences meet the performance criteria in at least one of the two columns except for the cases described in this document. Results for Endrin ketone qualified as estimated (J or UJ) in affected samples.

Several analytes missed the % difference continuing calibration criteria in cc1931-40 analyzed on 03/28/17 (9:49:54 PM). No samples from this batch associated with this continuing calibration check.

Final calibration verification included in data package.

Criteria

Are a five point calibration curve delivered with concentration levels as shown in Table 3 of SOP HW-36A, Revision 0, June, 2015? **Yes**? or No?

Actions

If the standard concentrations listed in Table 3 are not used, use professional judgment to evaluate the effect on the data

Criteria

Are RT Windows calculated correctly? **Yes**? or No?

Action

Recalculate the windows and use the corrected values for all evaluations.

Criteria

Are the Percent Relative Standard Deviation (%RSD) of the CFs for each of the single component target compounds less than or equal to 20.0%, except for alpha-BHC and delta-BHC?

Yes? or No?

All criteria were met __X__
Criteria were not met
and/or see below _____

Are the %RSD of the CFs for alpha-BHC and delta-BHC less than or equal to 25.0%. **Yes**? or No?

Is the %RSD of the CFs for each of the Toxaphene peaks must be < 30% when 5-point ICAL is performed? **Yes**? or No?

Is the %RSD of the CFs for the two surrogates (tetrachloro-m-xylene and decachlorobiphenyl) less than or equal to 30.0%. **Yes**? or No?

Action

a. If the %RSD criteria are not met, qualify detects as estimated (J) and use professional judgment to qualify non-detected target compounds.
b. If the %RSD criteria are within allowable limits, no qualification of the data is necessary

Continuing Calibration Checks

Criteria

Is the continuing calibration standard analyzed at the acceptable time intervals? **Yes**? or No?

Action

a. If more than 14 hours has elapsed from the injection of the instrument blank that begins an analytical sequence (opening CCV) and the injection of either a PEM or mid-point concentration of the Individual Standard Mixtures (A and B) or (C), qualify all data as unusable (R).
b. If more than 12 hours has elapsed from the injection of the instrument blank that begins an analytical sequence (opening CCV) and the injection of the last sample or blank that is part of the same analytical sequence, qualify all data as unusable (R).
c. If more than 72 hours has elapsed from the injection of the sample with a Toxaphene detection and the Toxaphene Calibration Verification Standard (CS3), qualify all data as unusable (R).

Criteria

Is the Percent Difference (%D) within ±25.0% for the PEM sample? Yes? or **No**?

Action
a. Qualify associated detects as estimated (J) and non-detects as estimated (UJ).

Criteria

For the Calibration Verification Standard (CS3); is the Percent Difference (%D) within ± 25.0%? Yes? or **No**?

Action

Qualify associated detects as estimated (J) and non-detects as estimated (UJ).

Criteria

Is the PEM 4,4'-DDT % Breakdown >20.0% and 4,4'-DDT is detected? Yes? or **No**?

Action

a. Qualify detects for 4,4'-DDT; detects for 4,4'-DDD; and detects for 4,4'-DDE as estimated (J)
b. Non-detected associated compounds are not qualified

Criteria

Is the PEM 4,4'-DDT % Breakdown >20.0% and 4,4'-DDT is not detected Yes? or **No**?

Action

a. Qualify non-detects for 4,4'- DDT as unusable (R)
b. Qualify detects for 4,4'-DDD as tentatively identified (NJ)
c. Qualify detects for 4,4'-DDE as tentatively identified (NJ)

Criteria

Is the PEM Endrin % Breakdown >20.0% and Endrin is detected? Yes? or **No**?

Action

a. Qualify detects for Endrin; detects for Endrin aldehyde; and detects for Endrin ketone as estimated (J)
b. Non-detected associated compounds are not qualified

Criteria

Is the PEM Endrin % Breakdown >20.0% and Endrin is not detected Yes? or **No**?

Action

a. Qualify non-detects for Endrin as unusable (R)
b. Qualify detects for Endrin aldehyde as tentatively identified (NJ)
c. Qualify detects for Endrin ketone as tentatively identified (NJ)

A separate worksheet should be filled for each initial curve

All criteria were met __X__
Criteria were not met
and/or see below _____

BLANK ANALYSIS RESULTS (Sections 1 & 2)

The assessment of the blank analysis results is to determine the existence and magnitude of contamination problems. The criteria for evaluation of blanks apply only to blanks associated with the samples, including trip, equipment, and laboratory blanks. If problems with any blanks exist, all data associated with the case must be carefully evaluated to determine whether or not there is an inherent variability in the data for the case, or if the problem is an isolated occurrence not affecting other data.

List the contamination in the blanks below. High and low levels blanks must be treated separately.

CRQL concentration_____N/A_____

Laboratory blanks

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
_No_target_analytes_detected_in_method_blanks_at_a_reporting_limit_of_0.01,_0.04,_and_0.25__				
_ug/L_____				

Field/Equipment/Trip blank

DATE ANALYZED	LAB ID	LEVEL/ MATRIX	COMPOUND	CONCENTRATION UNITS
_No_target_analytes_detected_in_the_field_blanks_analyzed_with_this_data_package._No_____				
_equipment_blank_analyzed_with_this_data_package._____				

All criteria were met __X__
Criteria were not met
and/or see below _____

BLANK ANALYSIS RESULTS (Section 3)

Blank Actions

Action Levels (ALs) should be based upon the highest concentration of contaminant determined in any blank. Do not qualify any blank with another blank. The ALs for samples which have been diluted should be corrected for the sample dilution factor and/or % moisture, where applicable. No positive sample results should be reported unless the concentration of the compound in the samples exceeds the ALs:

The concentration of non-target compounds in all blanks must be less than or equal to 10 µg/L. The concentration of each target compound found in the method or field blanks must be less than its CRQL listed in the method.

Data concerning the field blanks are not evaluated as part of the CCS process. If field blanks are present, the data reviewer should evaluate this data in a similar fashion as the method blanks.

Specific actions are as follows:

Blank Actions for Pesticide Analyses

Blank Type	Blank Result	Sample Result	Action for Samples
Method, Sulfur Cleanup, Instrument, Field, TCLP/SPLP	Detects	Not detected	No qualification required
	< CRQL	< CRQL	Report CRQL value with a U
		≥ CRQL	No qualification required
	> CRQL	< CRQL	Report CRQL value with a U
		≥ CRQL and ≤ blank concentration	Report blank value for sample concentration with a U
		≥ CRQL and > blank concentration	No qualification required
	= CRQL	≤ CRQL	Report CRQL value with a U
		> CRQL	No qualification required
	Gross contamination	Detects	Report blank value for sample concentration with a U

DATA REVIEW WORKSHEETS

All criteria were met __X__
Criteria were not met
and/or see below _____

CONTAMINATION SOURCE/LEVEL	COMPOUND	CONC/UNITS	AL/UNITS	SQL	AFFECTED SAMPLES

All criteria were met __X__
Criteria were not met
and/or see below _____

SURROGATE SPIKE RECOVERIES

Laboratory performance of individual samples is established by evaluation of surrogate spike recoveries. All samples are spiked with surrogate compounds prior to sample analysis. The accuracy of the analysis is measured by the surrogate percent recovery. Since the effects of the sample matrix are frequently outside the control of the laboratory and may present relatively unique problems, the validation of data is frequently subjective and demands analytical experience and professional judgment.

List the percent recoveries (%Rs) which do not meet the criteria for surrogate recovery.

Matrix:_Aqueous/Solid_____

Lab Sample ID	Lab File ID	S1 a	S2 a
FA42192-1	TT381599.D	111	121
FA42192-2	TT381600.D	106	130*
FA42192-3	TT381601.D	115	133*
FA42192-6	KK82632.D	98	113
FA42192-7	TT381607.D	110	127
FA42192-8	TT381608.D	115	115
FA42192-11	TT381609.D	107	102
FA42192-12	TT381610.D	132*	153*
FA42192-13	TT381611.D	116	135* b
FA42192-14	TT381612.D	115	129* b
FA42192-15	TT381613.D	133* b	120
OP64318-BS	TT381596.D	111	132*
OP64318-BS2	TT381597.D	111	129*
OP64318-MB	TT381598.D	108	135*
OP64318-MS	TT381602.D	107	135*
OP64318-MSD	TT381603.D	117	133*

Surrogate Compounds	Recovery Limits (Aqueous)
S1 = Tetrachloro-m-xylene	42-127%
S2 = Decachlorobiphenyl	27-127%

(a) Recovery from GC signal #1
(b) Outside control limits.

Note: Surrogate recoveries were within laboratory control limits except in the cases described in this document. No action taken, professional judgment. Surrogates recovered high and analytes not detected in affected samples.

Actions:

a. For any surrogate recovery greater than 150%, qualify detected target compounds as biased high (J+).
b. Do not qualify non-detected target compounds for surrogate recovery > 150 %.
c. If both surrogate recoveries are greater than or equal to 30% and less than or equal to 150%, no qualification of the data is necessary.
d. For any surrogate recovery greater than or equal to 10% and less than 30%, qualify detected target compounds as biased low (J-).
e. For any surrogate recovery greater than or equal to 10% and less than 30%, qualify non-detected target compounds as approximated (UJ).

f. If low surrogate recoveries are from sample dilution, professional judgment should be used to determine if the resulting data should be qualified. If sample dilution is not a factor:

i. Qualify detected target compounds as biased low (J-).
ii. Qualify non-detected target compounds as unusable (R).

g. If surrogate RTs in PEMs, Individual Standard Mixtures, samples, and blanks are outside of the RT Windows, the reviewer must use professional judgment to qualify data.
h. If surrogate RTs are within RT windows, no qualification of the data is necessary.
i. If the two surrogates were not added to all samples, MS/MSDs, standards, LCSs, and blanks, use professional judgment in qualifying data as missing surrogate analyte may not directly apply to target analytes.

Summary Surrogate Actions for Pesticide Analyses

Criteria	Action*	
	Detected Target Compounds	Non-detected Target Compounds
%R > 150%	J+	No qualification
30% < %R < 150%	No qualification	
10% < %R < 30%	J-	UJ
%R < 10% (sample dilution not a factor)	J-	R
%R < 10% (sample dilution is a factor)	Use professional judgment	
RT out of RT window	Use professional judgment	
RT within RT window	No qualification	

* Use professional judgment in qualifying data, as surrogate recovery problems may not directly apply to target analytes.

All criteria were met ___X___
Criteria were not met
and/or see below _____

MATRIX SPIKE/MATRIX SPIKE DUPLICATE (MS/MSD)

This data is generated to determine long term precision and accuracy in the analytical method for various matrices. This data alone cannot be used to evaluate the precision and accuracy of individual samples. If any % R in the MS or MSD falls outside the designated range, the reviewer should determine if there are matrix effects, i.e. LCS data are within the QC limits but MS/MSD data are outside QC limit.

1. MS/MSD Recoveries and Precision Criteria

Data for MS and MSDs will not be present unless requested by the Region.
Notify the Contract Laboratory Program Project Officer (CLP PO) if a field blank was used for the MS and MSD, unless designated as such by the Region.

> **NOTE:** For a Matrix Spike that does not meet criteria, apply the action to only the field sample used to prepare the Matrix Spike sample. If it is clearly stated in the data validation materials that the samples were taken through incremental sampling or some other method guaranteeing the homogeneity of the sample group, then the entire sample group may be qualified.

List the %Rs, RPD of the compounds which do not meet the criteria.

Sample ID:____FA42192-3MS/3MSD__ Matrix/Level:___Groundwater___

The QC reported here applies to the following samples: **Method:** SW846 8081B
FA42192-1, FA42192-2, FA42192-3, FA42192-6, FA42192-7, FA42192-8, FA42192-11, FA42192-12, FA42192-13 FA42192-14, FA42192-15

Compound	ug/l	Q	Spike ug/l	MS ug/l	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
Aldrin	ND		1	1.2	120	1	1.3	130*	8	61-126/21
alpha-BHC	ND		1	1.2	120	1	1.3	130*	8	66-129/23
alpha-Chlordane	ND		1	1.4	140*	1	1.5	150*	7	66-131/24
γ-Chlordane	0.017	J	1	1.7	167*	1	2.0	197*	16	68-128/21
Dieldrin	ND		1	1.3	130	1	1.4	140*	7	66-138/22
4,4'-DDD	ND		1	1.4	140*	1	1.5	150*	7	63-138/24
4,4'-DDE	ND		1	1.3	130	1	1.4	140*	7	59-133/23
4,4'-DDT	ND		1	1.4	140	1	1.5	150*	7	55-145/27
Endrin	ND		1	1.4	140	1	1.5	150*	7	71-147/23
Endosulfan sulfate	ND		1	1.4	140*	1	1.5	150*	7	64-128/26
Endrin ketone	ND		1	1.5	150*	1	1.5	150*	0	66-137/25
Endosulfan-II	ND		1	1.3	130	1	1.4	140*	7	65-133/22
Heptachlor epoxide	ND		1	1.2	120	1	1.3	130*	8	67-129/23
Methoxychlor	ND		1	1.4	140*	1	1.5	150*	7	60-136/25

Note: MS/MSD % recoveries and RPD within laboratory control limits except for the cases described in this document. No qualification made professional judgment. Analytes recovered high and not detected in sample batch above reporting limit except for alpha-Chlordane.

MS/MSD % recovery results apply to unspiked sample. Positive alpha-Chlordane results were from samples FA42192-7 and FA42192-12. No qualification made.

Action

No qualification of the data is necessary on MS and MSD data alone. However, using professional judgment, the validator may use the MS and MSD results in conjunction with other QC criteria and determine the need for some qualification of the data.

A separate worksheet should be used for each MS/MSD pair.

All criteria were met __X__
Criteria were not met
and/or see below _____

LABORATORY CONTROL SAMPLE (LCS) ANALYSIS

This data is generated to determine accuracy of the analytical method for various matrices.

1. LCS Recoveries Criteria

LCS Spike Compound	Recovery Limits (%)
gamma-BHC	50 – 120
Heptachlor epoxide	50 – 150
Dieldrin	30 – 130
4,4'-DDE	50 – 150
Endrin	50 – 120
Endosulfan sulfate	50 – 120
trans-Chlordane	30 – 130
Tetrachloro-m-xylene (surrogate)	30 – 150
Decachlorobiphenyl (surrogate)	30 – 150

LCS concentrations:____0.040_ug/l_____

List the %R of compounds which do not meet the criteria

LCS ID COMPOUND % R QC LIMIT

___%_recovery_and_RPD_within_laboratory_control_limits_except_for_the_cases_described_in___
___this_document._____
_____OP64318-BS___gamma-Chlordane_____160_%_____68_-_188_____
_____4,4'-DDD_____140_%_____63_-_138_____
_____Endosulfan_sulfate_____130_%_____64_-_128_____
_____Endosulfan_ketone_____140_%_____66_-_137_____

Note: LCS recovery exceeds the upper acceptance limit; non-detected compounds are not qualified.

Action

The following guidance is suggested for qualifying sample data for which the associated LCS does not meet the required criteria.

a. If the LCS recovery exceeds the upper acceptance limit, qualify detected target compounds as estimated (J). Do not qualify non-detected target compounds.

b. If the LCS recovery is less than the lower acceptance limit, qualify detected target compounds as estimated (J) and non-detects as unusable (R).

c. Use professional judgment to qualify data for compounds other than those compounds that are included in the LCS.

d. Use professional judgment to qualify non-LCS compounds. Take into account the compound class, compound recovery efficiency, analytical problems associated with each compound, and comparability in the performance of the LCS compound to the non-LCS compound.
e. If the LCS recovery is within allowable limits, no qualification of the data is necessary.

2. Frequency Criteria:

Where LCS analyzed at the required frequency and for each matrix? **Yes** or No.
If no, the data may be affected. Use professional judgment to determine the severity of the effect and qualify data accordingly. Discuss any actions below and list the samples affected.

All criteria were met _____
Criteria were not met
and/or see below __N/A_____

FLORISIL CARTRIDGE PERFORMANCE CHECK

NOTE: Florisil cartridge cleanup is mandatory for all extracts.

Criteria

Is the Florisil cartridge performance check conducted at least once on each lot of cartridges used for sample cleanup or every 6 months, whichever is most frequent? Yes? or No? **N/A**

Criteria

Are the results for the Florisil Cartridge Performance Check solution included with the data package? Yes? or No? **N/A**

Note: If % criteria are not met, examine the raw data for the presence of polar interferences and use professional judgment in qualifying the data as follows:

Action:

a. If the Percent Recovery is greater than 120% for any of the pesticide target compounds in the Florisil Cartridge Performance Check, qualify detected compounds as estimated (J). Do not qualify non-detected target compounds.
b. If the Percent Recovery is greater than or equal to 80% and less than or equal to 120% for all the pesticide target compounds, no qualification of the data is necessary.
c. If the Percent Recovery is greater than or equal to 10% and less than 80% for any of the pesticide target compounds in the Florisil Cartridge Performance Check, qualify detected target compounds as estimated (J) and non-detected target compounds as approximated (UJ).
d. If the Percent Recovery is less than 10% for any of the pesticide target compounds in the Florisil Cartridge Performance Check, qualify detected compounds as estimated (J) and qualify non-detected target compounds as unusable (R).
e. If the Percent Recovery of 2,4,5-trichlorophenol in the Florisil Cartridge Performance Check is greater than or equal to 5%, use professional judgment to qualify detected and non-detected target compounds, considering interference on the sample chromatogram.

Note: State in the Data Review Narrative potential effects on the sample data resulting from the Florisil Cartridge Performance Check analysis not yielding acceptable results.

Note:_ No information for Florisil cartridge performance check included in data package.

All criteria were met _____
Criteria were not met
and/or see below _____

GEL PERMEATION CHROMATOGRAPHY (GPC) PERFORMANCE CHECK

NOTE: GPC cleanup is mandatory for all soil samples.

If GPC criteria are not met, examine the raw data for the presence of high molecular weight contaminants; examine subsequent sample data for unusual peaks; and use professional judgment in qualifying the data. Notify the Contract Laboratory Program Project Officer (CLP PO) if the laboratory chooses to analyze samples under unacceptable GPC criteria.

Action:

a. If the Percent Recovery is less than 10% for the pesticide compounds and surrogates during the GPC calibration check, the non-detected target compounds may be suspect, qualify detected compounds as estimated (J).
b. If the Percent Recovery is less than 10% for the pesticide compounds and surrogates during the GPC calibration check, qualify all non-detected target compounds as unusable (R).
c. If the Percent Recovery is greater than or equal to 10% and is less than 80% for any of the pesticide target compounds in the GPC calibration, qualify detected target compounds as estimated (J) and non-detected target compounds as approximated (UJ).
d. If the Percent Recovery is greater than or equal to 80% and less than or equal to 120% for all the pesticide target compounds, no qualification of the data is necessary.
e. If high recoveries (i.e., greater than 120%) were obtained for the pesticides and surrogates during the GPC calibration check, qualify detected compounds as estimated (J). Do not qualify non-detected target compounds.

Note: State in the Data Review Narrative potential effects on the sample data resulting from the GPC cleanup analyses not yielding acceptable results.

Note:_ No information for performance of GPC cleanup included in data package.

All criteria were met __X__
Criteria were not met
and/or see below _____

TARGET COMPOUND IDENTIFICATION

Criteria:

1. Is Retention Times (RTs) of both of the surrogates and reported target compounds in each sample within the calculated RT Windows on both columns? **Yes**? or No?

2. Is the Tetrachloro-m-xylene (TCX) RT ±0.05 minutes of the Mean RT (RT) determined from the initial calibration and Decachlorobiphenyl (DCB) within ±0.10 minutes of the RT determined from the initial calibration? **Yes**? or No?

3. Is the Percent Difference (%D) for the detected mean concentrations of a pesticide target compound between the two Gas Chromatograph (GC) columns within the inclusive range of ± 25.0 %? **Yes**? or No?

4. When no analytes are identified in a sample; are the chromatograms from the analyses of the sample extract and the low-point standard of the initial calibration associated with those analyses on the same scaling factor? **Yes**? or No?

5. Does the chromatograms display the Single Component Pesticides (SCPs) detected in the sample and the largest peak of any multi-component analyte detected in the sample at less than full scale. Yes? or **No**?

6. If an extract is diluted; does the chromatogram display SCPs peaks between 10-100% of full scale, and multi-component analytes between 25-100% of full scale? Yes? or No? **N/A**

7. For any sample; does the baseline of the chromatogram return to below 50% of full scale before the elution time of alpha-BHC, and also return to below 25% of full scale after the elution time of alpha-BHC and before the elution time of DCB? **Yes**? or No?

8. If a chromatogram is replotted electronically to meet these requirements; is the scaling factor used displayed on the chromatogram, and both the initial chromatogram and the replotted chromatogram submitted in the data package. Yes? or **No**?

Action:

a. If the qualitative criteria for both columns were not met, all target compounds that are reported as detected should be considered non-detected.

b. Use professional judgment to assign an appropriate quantitation limit using the following guidance:

   i. If the detected target compound peak was sufficiently outside the pesticide RT Window, the reported values may be a false positive and should be replaced with the sample Contract Required Quantitation Limits (CRQL) value.

ii. If the detected target compound peak poses an interference with potential detection of another target peak, the reported value should be considered and qualified as unusable (R).

c. If the data reviewer identifies a peak in both GC column analyses that falls within the appropriate RT Windows, but was reported as a non-detect, the compound may be a false negative. Use professional judgment to decide if the compound should be included.

Note: State in the Data Review Narrative all conclusions made regarding target compound identification.

d. If the Toxaphene peak RT windows determined from the calibration overlap with SCPs or chromatographic interferences, use professional judgment to qualify the data.

e. If target compounds were detected on both GC columns, and the Percent Difference between the two results is greater than 25.0%, consider the potential for coelution and use professional judgment to decide whether a much larger concentration obtained on one column versus the other indicates the presence of an interfering compound. If an interfering compound is indicated, use professional judgment to determine how best to report, and if necessary, qualify the data according to these guidelines.

f. If Toxaphene exhibits a marginal pattern-matching quality, use professional judgment to establish whether the differences are due to environmental "weathering" (i.e., degradation of the earlier eluting peaks relative to the later eluting peaks). If the presence of Toxaphene is strongly suggested, report results as presumptively present (N).

GAS CHROMATOGRAPH/MASS SPECTROMETER (GC/MS) CONFIRMATION

NOTE: This confirmation is not usually provided by the laboratory. In cases where it is provided, use professional judgment to determine if data qualified with "C" can be salvaged if it was previously qualified as unusable (R).

Action:

a. If the quantitative criteria for both columns were met (≥ 5.0 ng/µL for SCPs and ≥ 125 ng/µL for Toxaphene), determine whether GC/MS confirmation was performed. If it was performed, qualify the data using the following guidance:

i. If GC/MS confirmation was not required because the quantitative criteria for both columns was not met, but it was still performed, use professional judgment when evaluating the data to decide whether the detect should be qualified with "C".

ii. If GC/MS confirmation was performed, but unsuccessful for a target compound detected by GC/ECD analysis, qualify those detects as "X".

All criteria were met __X__
Criteria were not met
and/or see below _____

COMPOUND QUANTITATION AND REPORTED CONTRACT REQUIRED QUANTITATION LIMITS (CRQLS)

The sample quantitation evaluation is to verify laboratory quantitation results. In the space below, please show a minimum of one sample calculation:

FA42192-7 alpha-chlordane RF = 5.554×10^4

[ ] = $(98591)/(5.554 \times 10^4)$
= 1.78 ppb Ok

Note:

Action:

a. If sample quantitation is different from the reported value, qualify result as unusable (R).
b. When a sample is analyzed at more than one dilution, the lowest CRQLs are used unless a QC exceedance dictates the use of the higher CRQLs from the diluted sample.
c. Replace concentrations that exceed the calibration range in the original analysis by crossing out the "E" and its corresponding value on the original reporting form and substituting the data from the diluted sample.
d. Results between the MDL and CRQL should be qualified as estimated (J).
e. Results less than the MDL should be reported at the CRQL and qualified (U). MDLs themselves are not reported.
f. For non-aqueous samples, if the percent moisture is less than 70.0%, no qualification of the data is necessary. If the percent moisture is greater than or equal to 70.0% and less than 90.0%, qualify detects as estimated (J) and non-detects as approximated (UJ). If the percent moisture is greater than or equal to 90.0%, qualify detects as estimated (J) and non-detects as unusable (R) (see Table).

Percent Moisture Actions for Pesticide Analysis for Non-Aqueous Samples

Criteria	**Action**	
	Detected Associated Compounds	**Non-detected Associated Compounds**
% Moisture < 70.0	No qualification	
70.0 < % Moisture < 90.0	J	UJ
% Moisture > 90.0	J	R

List samples which have ≤ 50 % solids

Note: If any discrepancies are found, the Region's designated representative may contact the laboratory to obtain additional information that could resolve any differences. If a discrepancy remains unresolved, the reviewer must use professional judgment to decide which value is the most accurate. Under these circumstances, the reviewer may determine that qualification of data is warranted. Note in the Data Review Narrative a description of the reasons for data qualification and the qualification that is applied to the data.

Dilution performed

SAMPLE ID	DILUTION FACTOR	REASON FOR DILUTION

All criteria were met __X_____
Criteria were not met
and/or see below _____

FIELD DUPLICATE PRECISION

NOTE: In the absence of QAPP guidance for validating data from field duplicates, the following action will be taken.

Field duplicates samples may be taken and analyzed as an indication of overall precision. These analyses measure both field and lab precision; therefore, the results may have more variability than laboratory duplicates which only laboratory performance. It is also expected that soil duplicate results will have a greater variance than water matrices due to difficulties associated with collecting identical field duplicate samples. Identify which samples within the data package are field duplicates. Estimate the relative percent difference (RPD) between the values for each compound. If large RPDs (> 50%) is observed, confirm identification of samples and note difference in the executive summary.

Sample IDs: _FA42192-8/FA42192-11_____ Matrix:__Groundwater__

COMPOUND	SQL ug/L	SAMPLE CONC.	DUPLICATE CONC.	RPD	ACTION
Field duplicate analyzed with this data package. RPD within the required criteria of < 50 %.					

Actions:

a. Qualify as estimated positive results (J) and nondetects (UJ) for the compound that exceeded the above criteria. For organics, only the sample and duplicate will be qualified.

b. If an RPD cannot be calculated because one or both of the sample results is not detected, the following actions apply:

i. If one sample result is not detected and the other is greater than 5x the SQL qualify (J/UJ).

ii. If one sample value is not detected and the other is greater than 5x the SQL and the SQLs for the sample and duplicate are significantly different, use professional judgment to determine if qualification is appropriate.

iii. If one sample value is not detected and the other is less than 5x, use professional judgment to determine if qualification is appropriate.

iv. If both sample and duplicate results are not detected, no action is needed.

OVERALL ASSESSMENT OF DATA
Action:

1. Use professional judgment to determine if there is any need to qualify data which were not qualified based on the Quality Control (QC) criteria previously discussed.

2. Write a brief narrative to give the user an indication of the analytical limitations of the data.

Note: The Contract Laboratory Program Project Officer (CLP PO) must be informed if any inconsistency of the data with the Sample Delivery Group (SDG) Narrative. If sufficient information on the intended use and required quality of the data is available, the reviewer should include their assessment of the usability of the data within the given context. This may be used as part of a formal Data Quality Assessment (DQA).

Overall assessment of the data: Results are valid; the data can be used for decision making purposes.