



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:52:02PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-26-2

Laboratory: BC Laboratories SDG: 17-20817
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720817-07 File ID: 02AUG49.D
Sampled: 07/28/17 10:50 Prepared: 08/02/17 07:00 Analyzed: 08/03/17 04:02
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0012 Sequence: 1713601 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	1.3	
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

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3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:52:02PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-26-2

Laboratory: BC Laboratories SDG: 17-20817
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720817-07 File ID: 02AUG49.D
Sampled: 07/28/17 10:50 Prepared: 08/02/17 07:00 Analyzed: 08/03/17 04:02
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0012 Sequence: 1713601 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	1.8	
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.28	J
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

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Project Number: 3Q17
Project Manager: David Conner

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MW-26-2

Laboratory: BC Laboratories SDG: 17-20817
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720817-07 File ID: 02AUG49.D
Sampled: 07/28/17 10:50 Prepared: 08/02/17 07:00 Analyzed: 08/03/17 04:02
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0012 Sequence: 1713601 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U UT
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U UT
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.580	106	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.7700	97.7	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.6000	96.0	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	193818	6.57	204288	6.57	
Chlorobenzene-d5 (IS)	74170	9.61	82858	9.61	
1,4-Difluorobenzene (IS)	276070	7.38	309776	7.38	

* Values outside of QC limits

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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-26-2

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-20817</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1720817-07</u>	File ID:	<u>02AUG49.D</u>		
Sampled:	<u>07/28/17 10:50</u>	Prepared:	<u>08/02/17 07:00</u>	Analyzed:	<u>08/03/17 04:02</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0012</u>	Sequence:	<u>1713601</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloro propanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

SL 9/2/17



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ORGANIC ANALYSIS DATA SHEET
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MW-26-1

Laboratory: BC Laboratories SDG: 17-20817
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720817-08 File ID: 02AUG50.D
Sampled: 07/28/17 11:10 Prepared: 08/02/17 07:00 Analyzed: 08/03/17 04:25
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0012 Sequence: 1713601 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.29	J
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:52:02PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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EPA-524.2

MW-26-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-08</u>	File ID: <u>02AUG50.D</u>	
Sampled: <u>07/28/17 11:10</u>	Prepared: <u>08/02/17 07:00</u>	Analyzed: <u>08/03/17 04:25</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BJH0012</u>	Sequence: <u>1713601</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.70	
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.24	J
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

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Tidewater Inc. Reported: 8/25/2017 12:52:02PM
 3761 Attucks Drive Project: JPL- GW Monitoring Wells
 Powell, OH 43065 Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-26-1

Laboratory: BC Laboratories SDG: 17-20817
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720817-08 File ID: 02AUG50.D
 Sampled: 07/28/17 11:10 Prepared: 08/02/17 07:00 Analyzed: 08/03/17 04:25
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0012 Sequence: 1713601 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U UJ
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U UJ
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	11.290	113	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.7700	97.7	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	10.200	102	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	183962	6.57	204288	6.57	
Chlorobenzene-d5 (IS)	71799	9.61	82858	9.61	
1,4-Difluorobenzene (IS)	271114	7.38	309776	7.38	

* Values outside of QC limits

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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-26-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-08</u>	File ID: <u>02AUG50.D</u>	
Sampled: <u>07/28/17 11:10</u>	Prepared: <u>08/02/17 07:00</u>	Analyzed: <u>08/03/17 04:25</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BIH0012</u>	Sequence: <u>1713601</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

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Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-5-072817

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-20817</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1720817-09</u>	File ID:	<u>02AUG51.D</u>		
Sampled:	<u>07/28/17 11:25</u>	Prepared:	<u>08/02/17 07:00</u>	Analyzed:	<u>08/03/17 04:48</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0012</u>	Sequence:	<u>1713601</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/25/17



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Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
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EB-5-072817

Laboratory: BC Laboratories SDG: 17-20817
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720817-09 File ID: 02AUG51.D
Sampled: 07/28/17 11:25 Prepared: 08/02/17 07:00 Analyzed: 08/03/17 04:48
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0012 Sequence: 1713601 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,1,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

SL 89217



Tidewater Inc. Reported: 8/25/2017 12:52:02PM
 3761 Attucks Drive Project: JPL- GW Monitoring Wells
 Powell, OH 43065 Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-5-072817

Laboratory: BC Laboratories SDG: 17-20817
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720817-09 File ID: 02AUG51.D
 Sampled: 07/28/17 11:25 Prepared: 08/02/17 07:00 Analyzed: 08/03/17 04:48
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0012 Sequence: 1713601 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UT</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UT</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	11.340	113	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.9600	99.6	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	10.230	102	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	183371	6.57	204288	6.57	
Chlorobenzene-d5 (IS)	73943	9.61	82858	9.61	
1,4-Difluorobenzene (IS)	276989	7.38	309776	7.38	

* Values outside of QC limits

SL09217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:52:02PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-5-072817

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-09</u>	File ID: <u>02AUG51.D</u>	
Sampled: <u>07/28/17 11:25</u>	Prepared: <u>08/02/17 07:00</u>	Analyzed: <u>08/03/17 04:48</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BIH0012</u>	Sequence: <u>1713601</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8292117

LDC #: 39385A1

VALIDATION COMPLETENESS WORKSHEET

SDG #: 17-20817

Level III

Laboratory: BC Laboratories, Inc.

Date: 9/12/17

Page: 1 of 1

Reviewer: [Signature]

2nd Reviewer: [Signature]

METHOD: GC/MS Volatiles (EPA Method 524.2)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A	
II.	GC/MS Instrument performance check	A	
III.	Initial calibration/ICV	A/A	RSD ≤ 20%. Y ² ICV = 35%
IV.	Continuing calibration	MW	CV = 35%
V.	Laboratory Blanks	A	
VI.	Field blanks	ND	TB=1. EB=8
VII.	Surrogate spikes	A	
VIII.	Matrix spike/Matrix spike duplicates	N	CS
IX.	Laboratory control samples	A	LCS
X.	Field duplicates	MW	D = 3+5
XI.	Internal standards	A	
XII.	Compound quantitation RL/LOQ/LODs	N	
XIII.	Target compound identification	N	
XIV.	System performance	N	
XV.	Overall assessment of data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	TB-5-072817	1720817-01	Water	07/28/17
2	MW-23-3	1720817-03	Water	07/28/17
3	MW-23-2	1720817-04	Water	07/28/17
4	MW-23-1	1720817-05	Water	07/28/17
5	DUPE -3-3Q17	1720817-06	Water	07/28/17
6	MW-26-2	1720817-07	Water	07/28/17
7	MW-26-1	1720817-08	Water	07/28/17
8	EB-5-072817	1720817-09	Water	07/28/17
9				
10				

Notes:

TARGET COMPOUND WORKSHEET

METHOD: VOA

A. Chloromethane	AA. Tetrachloroethene	AAA. 1,3,5-Trimethylbenzene	AAAA. Ethyl tert-butyl ether	A1. 1,3-Butadiene
B. Bromomethane	BB. 1,1,2,2-Tetrachloroethane	BBB. 4-Chlorotoluene	BBBB. tert-Amyl methyl ether	B1. Hexane
C. Vinyl chloride	CC. Toluene	CCC. tert-Butylbenzene	CCCC. 1-Chlorohexane	C1. Heptane
D. Chloroethane	DD. Chlorobenzene	DDD. 1,2,4-Trimethylbenzene	DDDD. Isopropyl alcohol	D1. Propylene
E. Methylene chloride	EE. Ethylbenzene	EEE. sec-Butylbenzene	EEEE. Acetonitrile	E1. Freon 11
F. Acetone	FF. Styrene	FFF. 1,3-Dichlorobenzene	FFFF. Acrolein	F1. Freon 12
G. Carbon disulfide	GG. Xylenes, total	GGG. p-Isopropyltoluene	GGGG. Acrylonitrile	G1. Freon 113
H. 1,1-Dichloroethene	HH. Vinyl acetate	HHH. 1,4-Dichlorobenzene	HHHH. 1,4-Dioxane	H1. Freon 114
I. 1,1-Dichloroethane	II. 2-Chloroethylvinyl ether	III. n-Butylbenzene	IIII. Isobutyl alcohol	I1. 2-Nitropropane
J. 1,2-Dichloroethene, total	JJ. Dichlorodifluoromethane	JJJ. 1,2-Dichlorobenzene	JJJJ. Methacrylonitrile	J1. Dimethyl disulfide
K. Chloroform	KK. Trichlorofluoromethane	KKK. 1,2,4-Trichlorobenzene	KKKK. Propionitrile	K1. 2,3-Dimethyl pentane
L. 1,2-Dichloroethane	LL. Methyl-tert-butyl ether	LLL. Hexachlorobutadiene	LLLL. Ethyl ether	L1. 2,4-Dimethyl pentane
M. 2-Butanone	MM. 1,2-Dibromo-3-chloropropane	MMM. Naphthalene	MMMM. Benzyl chloride	M1. 3,3-Dimethyl pentane
N. 1,1,1-Trichloroethane	NN. Methyl ethyl ketone	NNN. 1,2,3-Trichlorobenzene	NNNN. Iodomethane	N1. 2-Methylpentane
O. Carbon tetrachloride	OO. 2,2-Dichloropropane	OOO. 1,3,5-Trichlorobenzene	OOOO. 1,1-Difluoroethane	O1. 3-Methylpentane
P. Bromodichloromethane	PP. Bromochloromethane	PPP. trans-1,2-Dichloroethene	PPPP. Tetrahydrofuran	P1. 3-Ethylpentane
Q. 1,2-Dichloropropane	QQ. 1,1-Dichloropropene	QQQ. cis-1,2-Dichloroethene	QQQQ. Methyl acetate	Q1. 2,2-Dimethylpentane
R. cis-1,3-Dichloropropene	RR. Dibromomethane	RRR. m,p-Xylenes	RRRR. Ethyl acetate	R1. 2,2,3-Trimethylbutane
S. Trichloroethene	SS. 1,3-Dichloropropane	SSS. o-Xylene	SSSS. Cyclohexane	S1. 2,2,4-Trimethylpentane
T. Dibromochloromethane	TT. 1,2-Dibromoethane	TTT. 1,1,2-Trichloro-1,2,2-trifluoroethane	TTTT. Methylcyclohexane	T1. 2-Methylhexane
U. 1,1,2-Trichloroethane	UU. 1,1,1,2-Tetrachloroethane	UUU. 1,2-Dichlorotetrafluoroethane	UUUU. Allyl chloride	U1. Nonanal
V. Benzene	VV. Isopropylbenzene	VVV. 4-Ethyltoluene	VVVV. Methyl methacrylate	V1. 2-Methylnaphthalene
W. trans-1,3-Dichloropropene	WW. Bromobenzene	WWW. Ethanol	WWWW. Ethyl methacrylate	W1. Methanol
X. Bromoform	XX. 1,2,3-Trichloropropane	XXX. Di-isopropyl ether	XXXX. cis-1,4-Dichloro-2-butene	X1. 1,2,3-Trimethylbenzene
Y. 4-Methyl-2-pentanone	YY. n-Propylbenzene	YYY. tert-Butanol	YYYY. trans-1,4-Dichloro-2-butene	Y1.
Z. 2-Hexanone	ZZ. 2-Chlorotoluene	ZZZ. tert-Butyl alcohol	ZZZZ. Pentachloroethane	Z1.

**VALIDATION FINDINGS WORKSHEET
Continuing Calibration**

METHOD: GC/MS VOA (EPA Method 524.2)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

N/A Was a continuing calibration standard analyzed at least once every 12 hours for each instrument?

N/A Were all percent differences (%D) \leq 30% ?

#	Date	Standard ID	Compound	Finding %D (Limit: \leq 30.0%)	Associated Samples	Qualifications
	8/1/17	01AUG33	B Methyl iodide 2222	63.5 62.9 59.6	1, 2, MB (ND)	✓N/A ↓
	8/2/17	02AUG02	B Methyl iodide	68.5 73.9	BAK	✓N/A ↓
	8/2/17	02AUG34	Methyl iodide 2222	53.9 35.3	3-8	✓N/A ↓

LDC#: 39385A1

VALIDATION FINDINGS WORKSHEET
Field Duplicates

Page: 1 of 1
Reviewer: g
2nd Reviewer: AE

METHOD: GCMS VOA (EPA Method 524.2)

Compound	Concentration (ug/L)		RPD
	3	5	
K	0.84	0.89	6
I	0.30	0.31	3
AA	0.93	0.80	15
S	4.0	4.1	2

V:\FIELD DUPLICATES\Field Duplicates\FD_Organics\2017\39385A1_JPL.wpd

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017

LDC Report Date: September 21, 2017

Parameters: Chromium

Validation Level: Level III

Laboratory: BC Laboratories, Inc.

Sample Delivery Group (SDG): 17-20817

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-23-4	1720817-02	Water	07/28/17
MW-23-3	1720817-03	Water	07/28/17
MW-23-2	1720817-04	Water	07/28/17
MW-23-1	1720817-05	Water	07/28/17
DUP-3-3Q17	1720817-06	Water	07/28/17
MW-26-2	1720817-07	Water	07/28/17
MW-26-1	1720817-08	Water	07/28/17
EB-5-072817	1720817-09	Water	07/28/17
MW-23-4MS	1720817-02MS	Water	07/28/17
MW-23-4MSD	1720817-02MSD	Water	07/28/17
MW-23-4DUP	1720817-02DUP	Water	07/28/17

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following method:

Chromium by Environmental Protection Agency (EPA) Method 200.8

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. ICPMS Tune

The mass calibration was within 0.1 AMU and the percent relative standard deviation (%RSD) was less than or equal to 5%.

III. Instrument Calibration

Initial and continuing calibrations were performed as required by the method.

The initial calibration verification (ICV) and continuing calibration verification (CCV) standards were within QC limits.

IV. ICP Interference Check Sample Analysis

ICP interference check sample analyses were not required by the method.

V. Laboratory Blanks

Laboratory blanks were analyzed as required by the methods. No contaminants were found in the laboratory blanks.

VI. Field Blanks

Sample EB-5-072817 was identified as an equipment blank. No contaminants were found with the following exceptions:

Blank ID	Analyte	Concentration
EB-5-072817	Chromium	1.2 ug/L

VII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VIII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

IX. Serial Dilution

Serial dilution was not performed for this SDG.

X. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the method. Percent recoveries (%R) were within QC limits.

XI. Field Duplicates

Samples MW-23-2 and DUP-3-3Q17 were identified as field duplicates. No results were detected in any of the samples with the following exceptions:

Analyte	Concentration (ug/L)		RPD
	MW-23-2	DUP-3-3Q17	
Chromium	2.2	2.2	0

XII. Internal Standards (ICP-MS)

Raw data were not reviewed for Level III validation.

XIII. Sample Result Verification

Raw data were not reviewed for Level III validation.

XIV. Overall Assessment of Data

The analysis was conducted within all specifications of the method. No results were rejected in this SDG.

The quality control criteria reviewed were met and are considered acceptable. Based upon the data validation all results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Chromium - Data Qualification Summary - SDG 17-20817

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017
Chromium - Laboratory Blank Data Qualification Summary - SDG 17-20817

No Sample Data Qualified in this SDG



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 2:24:31PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-23-4

Laboratory: BC Laboratories

SDG: 17-20817

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720817-02

File ID: PE-EL3 170803R-111

Sampled: 07/28/17 07:25

Prepared: 08/03/17 08:30

Analyzed: 08/03/17 23:13

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0270

Sequence: 1713864

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	3.7	1		EPA-200.8

209217



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 2:24:31PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-23-3

Laboratory: BC Laboratories

SDG: 17-20817

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720817-03

File ID: PE-EL3_170803R-121

Sampled: 07/28/17 07:50

Prepared: 08/03/17 08:30

Analyzed: 08/03/17 23:47

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0270

Sequence:

1713864

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	3.8	1		EPA-200.8

8092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 2:24:31PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-23-2

Laboratory: BC Laboratories

SDG: 17-20817

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720817-04

File ID: PE-EL3_170803R-122

Sampled: 07/28/17 08:25

Prepared: 08/03/17 08:30

Analyzed: 08/03/17 23:50

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0270

Sequence:

1713864

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.2	1	J	EPA-200.8

08/25/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 2:24:31PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-23-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-05</u>	File ID: <u>PE-EL3 170803R-123</u>	
Sampled: <u>07/28/17 09:10</u>	Prepared: <u>08/03/17 08:30</u>	Analyzed: <u>08/03/17 23:54</u>	
Solids: <u>0.00</u>	Preparation: <u>EPA 200.2</u>	Initial/Final: <u>50 ml / 50 ml</u>	
Batch: <u>BIH0270</u>	Sequence: <u>1713864</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>PE-EL3</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.0	1	J	EPA-200.8

209-117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 2:24:31PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-200.8

DUP-3-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-06</u>	File ID: <u>PE-EL3 170803R-124</u>	
Sampled: <u>07/28/17 08:35</u>	Prepared: <u>08/03/17 08:30</u>	Analyzed: <u>08/03/17 23:57</u>	
Solids: <u>0.00</u>	Preparation: <u>EPA 200.2</u>	Initial/Final: <u>50 ml / 50 ml</u>	
Batch: <u>BIH0270</u>	Sequence: <u>1713864</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>PE-EL3</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.2	1	J	EPA-200.8

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 2:24:31PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-26-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-07</u>	File ID: <u>PE-EL3 170803R-125</u>	
Sampled: <u>07/28/17 10:50</u>	Prepared: <u>08/03/17 08:30</u>	Analyzed: <u>08/04/17 00:01</u>	
Solids: <u>0.00</u>	Preparation: <u>EPA 200.2</u>	Initial/Final: <u>50 ml / 50 ml</u>	
Batch: <u>BJH0270</u>	Sequence: <u>1713864</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>PE-EL3</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	3.0	1		EPA-200.8

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 2:24:31PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-26-1

Laboratory: BC Laboratories

SDG: 17-20817

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720817-08

File ID: PE-EL3 170803R-126

Sampled: 07/28/17 11:10

Prepared: 08/03/17 08:30

Analyzed: 08/04/17 00:04

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0270

Sequence:

1713864

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.6	1	J	EPA-200.8

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 2:24:31PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-200.8

EB-5-072817

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-09</u>	File ID: <u>PE-EL3_170803R-127</u>	
Sampled: <u>07/28/17 11:25</u>	Prepared: <u>08/03/17 08:30</u>	Analyzed: <u>08/04/17 00:08</u>	
Solids: <u>0.00</u>	Preparation: <u>EPA 200.2</u>	Initial/Final: <u>50 ml / 50 ml</u>	
Batch: <u>B[H0270]</u>	Sequence: <u>1713864</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>PE-EL3</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.2	1	J	EPA-200.8

8/29/17

LDC #: 39385A4a

VALIDATION COMPLETENESS WORKSHEET

Date: 9/18/17

SDG #: 17-20817

Level III

Page: 1 of 1

Laboratory: BC Laboratories, Inc.

Reviewer: JB

2nd Reviewer: KR

METHOD: Chromium (EPA Method 200.8)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A/A	
II.	ICP/MS Tune	A	
III.	Instrument Calibration	A	
IV.	ICP Interference Check Sample (ICS) Analysis	N	Not Required
V.	Laboratory Blanks	A	
VI.	Field Blanks	SW	EB = 8
VII.	Matrix Spike/Matrix Spike Duplicates	A	
VIII.	Duplicate sample analysis	A	
IX.	Serial Dilution	N	Not performed
X.	Laboratory control samples	A	LCS
XI.	Field Duplicates	SW	(3,5)
XII.	Internal Standard (ICP-MS)	N	Not Reviewed for Level III
XIII.	Sample Result Verification	N	
XIV.	Overall Assessment of Data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	MW-23-4	1720817-02	Water	07/28/17
2	MW-23-3	1720817-03	Water	07/28/17
3	MW-23-2	1720817-04	Water	07/28/17
4	MW-23-1	1720817-05	Water	07/28/17
5	DUP/3-3Q17	1720817-06	Water	07/28/17
6	MW-26-2	1720817-07	Water	07/28/17
7	MW-26-1	1720817-08	Water	07/28/17
8	EB-5-072817	1720817-09	Water	07/28/17
9	MW-23-4MS	1720817-02MS	Water	07/28/17
10	MW-23-4MSD	1720817-02MSD	Water	07/28/17
11	MW-23-4DUP	1720817-02DUP	Water	07/28/17
12				
13				
14				

Notes:

LDC #: _____
SDG #: _____

VALIDATION FINDINGS WORKSHEET

Field Blanks

Page: 1 of 1
Reviewer: YB
2nd reviewer: KK

METHOD: Trace Metals (EPA CLP SOW ILM02.1)

Y N N/A Were field blanks identified in this SDG?
Y N N/A Were target analytes detected in the field blanks?

Sample: 8 Field Blank / Trip Blank / Rinsate / Other EB (circle one)

Analyte	Concentration Units (<u>µg/L</u>)
<u>Cr</u>	<u>1.2</u>

Sample: _____ Field Blank / Trip Blank / Rinsate / Other _____ (circle one)

Analyte	Concentration Units ()

LDC#: 39385A4a

VALIDATION FINDINGS WORKSHEET
Field Duplicates

Page: 1 of 1
Reviewer: JB
2nd Reviewer: KK

METHOD: Metals (EPA Method 200.8)

Analyte	Concentration (ug/L)		RPD	
	3	5		
Chromium	2.2	2.2	0	

V:\FIELD DUPLICATES\Field Duplicates\FD_inorganic\2017\39385A4a.wpd

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017

LDC Report Date: September 21, 2017

Parameters: Wet Chemistry

Validation Level: Level III

Laboratory: BC Laboratories, Inc.

Sample Delivery Group (SDG): 17-20817

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-23-4	1720817-02	Water	07/28/17
MW-23-3	1720817-03	Water	07/28/17
MW-23-2	1720817-04	Water	07/28/17
MW-23-1	1720817-05	Water	07/28/17
DUP-3-3Q17	1720817-06	Water	07/28/17
MW-26-2	1720817-07	Water	07/28/17
MW-26-1	1720817-08	Water	07/28/17
EB-5-072817	1720817-09	Water	07/28/17
MW-23-4MS	1720817-02MS	Water	07/28/17
MW-23-4MSD	1720817-02MSD	Water	07/28/17
MW-23-4DUP	1720817-02DUP	Water	07/28/17

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following methods:

Hexavalent Chromium by Environmental Protection Agency (EPA) SW 846 Method 7196

Perchlorate by EPA Method 314.0

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. Initial Calibration

All criteria for the initial calibration of each method were met.

III. Continuing Calibration

Continuing calibration frequency and analysis criteria were met for each method when applicable.

IV. Laboratory Blanks

Laboratory blanks were analyzed as required by the methods. No contaminants were found in the laboratory blanks.

V. Field Blanks

Sample EB-5-072817 was identified as an equipment blank. No contaminants were found.

VI. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

VIII. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the methods. Percent recoveries (%R) were within QC limits.

IX. Field Duplicates

Samples MW-23-2 and DUPE-3-3Q17 were identified as field duplicates. No results were detected in any of the samples with the following exceptions:

Analyte	Concentration (mg/L)		RPD
	MW-23-2	DUP-3-3Q17	
Perchlorate	3.9 ug/L	4.7 ug/L	19
Hexavalent Chromium	0.0011 mg/L	0.00099 mg/L	11

X. Sample Result Verification

Raw data were not reviewed for Level III validation.

XI. Overall Assessment of Data

The analysis was conducted within all specifications of the methods. No results were rejected in this SDG.

The quality control criteria reviewed were met and are considered acceptable. Based upon the data validation all results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Wet Chemistry - Data Qualification Summary - SDG 17-20817

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 17-20817

No Sample Data Qualified in this SDG



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:54:01PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-314.0

MW-23-3

Laboratory: BC Laboratories

SDG: 17-20817

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720817-03

File ID: F080917.seq-8.0000.txt

Sampled: 07/28/17 07:50

Prepared: 08/09/17 12:00

Analyzed: 08/09/17 22:23

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H0982

Sequence: 1714391

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.7	1	J	EPA-314.0

17292117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-23-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-04RE1</u>	File ID: <u>F080917.seq-10.0000.txt</u>	
Sampled: <u>07/28/17 08:25</u>	Prepared: <u>08/09/17 12:00</u>	Analyzed: <u>08/09/17 22:50</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BJH0982</u>	Sequence: <u>1714391</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	3.9	1	J	EPA-314.0

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-314.0

MW-23-1

Laboratory: BC Laboratories

SDG: 17-20817

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720817-05

File ID: F080917.seq-57.0000.txt

Sampled: 07/28/17 09:10

Prepared: 08/09/17 12:00

Analyzed: 08/10/17 10:36

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BFH0982

Sequence: 1714391

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	3.9	1	J	EPA-314.0

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

DUP-3-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-06</u>	File ID: <u>F080917.seq-59.0000.txt</u>	
Sampled: <u>07/28/17 08:35</u>	Prepared: <u>08/09/17 12:00</u>	Analyzed: <u>08/10/17 11:03</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B[H0982</u>	Sequence: <u>1714391</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	4.7	1		EPA-314.0

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-26-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-07</u>	File ID: <u>F080917.seq-60.0000.txt</u>	
Sampled: <u>07/28/17 10:50</u>	Prepared: <u>08/09/17 12:00</u>	Analyzed: <u>08/10/17 11:17</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B[H0982]</u>	Sequence: <u>1714391</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	3.9	1	J	EPA-314.0

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-26-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-08</u>
Sampled: <u>07/28/17 11:10</u>	Prepared: <u>08/09/17 12:00</u>
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>
Batch: <u>BIH0982</u>	Sequence: <u>1714391</u>
	Calibration: <u>UNASSIGNED</u>
	Instrument: <u>IC6</u>
	File ID: <u>F080917.seq-62.0000.txt</u>
	Analyzed: <u>08/10/17 11:45</u>
	Initial/Final: <u>20 ml / 20 ml</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.4	1	J	EPA-314.0

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

EB-5-072817

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-09</u>	File ID: <u>F080917.seq-64.0000.txt</u>	
Sampled: <u>07/28/17 11:25</u>	Prepared: <u>08/09/17 12:00</u>	Analyzed: <u>08/10/17 12:13</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B[H0982]</u>	Sequence: <u>1714391</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

Handwritten: 8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-7196

MW-23-4

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-02</u>
Sampled: <u>07/28/17 07:25</u>	Prepared: <u>07/28/17 17:28</u>
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>
Batch: <u>BIG2617</u>	Sequence: <u>1713607</u>
	Calibration: <u>UNASSIGNED</u>
	Instrument: <u>KONE-1</u>
	File ID: <u>170728 1728 CR6-017</u>
	Analyzed: <u>07/28/17 18:25</u>
	Initial/Final: <u>20 ml / 20 ml</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0031	1		EPA-7196

8092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-7196

MW-23-3

Laboratory: BC Laboratories

SDG: 17-20817

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720817-03

File ID: 170728 1728 CR6-025

Sampled: 07/28/17 07:50

Prepared: 07/28/17 17:28

Analyzed: 07/28/17 18:39

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIG2617

Sequence: 1713607

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0029	1		EPA-7196

Handwritten: X-292117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-7196

MW-23-2

Laboratory: BC Laboratories

SDG: 17-20817

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720817-04

File ID: 170728 1728 CR6-044

Sampled: 07/28/17 08:25

Prepared: 07/28/17 17:28

Analyzed: 07/28/17 19:12

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BJG2617

Sequence: 1713607

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0011	1	J	EPA-7196

8092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:54:01PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-23-1

Laboratory: BC Laboratories

SDG: 17-20817

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720817-05

File ID: 170728 1728 CR6-027

Sampled: 07/28/17 09:10

Prepared: 07/28/17 17:28

Analyzed: 07/28/17 18:39

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIG2617

Sequence: 1713607

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

DUP-3-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-06</u>	File ID: <u>170728 1728 CR6-028</u>	
Sampled: <u>07/28/17 08:35</u>	Prepared: <u>07/28/17 17:28</u>	Analyzed: <u>07/28/17 18:39</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIG2617</u>	Sequence: <u>1713607</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00099	1	J	EPA-7196

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-26-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-07</u>	File ID: <u>170728 1728 CR6-029</u>	
Sampled: <u>07/28/17 10:50</u>	Prepared: <u>07/28/17 17:28</u>	Analyzed: <u>07/28/17 18:39</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIG2617</u>	Sequence: <u>1713607</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0013	1	J	EPA-7196

D69-117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-26-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-08</u>	File ID: <u>170728 1728 CR6-030</u>	
Sampled: <u>07/28/17 11:10</u>	Prepared: <u>07/28/17 17:28</u>	Analyzed: <u>07/28/17 18:39</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIG2617</u>	Sequence: <u>1713607</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:54:01PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

EB-5-072817

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20817</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720817-09</u>	File ID: <u>170728 1728 CR6-031</u>	
Sampled: <u>07/28/17 11:25</u>	Prepared: <u>07/28/17 17:28</u>	Analyzed: <u>07/28/17 18:39</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIG2617</u>	Sequence: <u>1713607</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

SL292117

LDC #: 39385A6

VALIDATION COMPLETENESS WORKSHEET

Date: 9/18/17

SDG #: 17-20817

Level III

Page: 1 of 1

Laboratory: BC Laboratories, Inc.

Reviewer: JB

2nd Reviewer: KE

METHOD: (Analyte) Hexavalent Chromium (EPA SW846 Method 7196), Perchlorate (EPA Method 314.0)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A / A	
II	Initial calibration	A	
III.	Calibration verification	A	
IV	Laboratory Blanks	A	
V	Field blanks	ND	EB = 8
VI.	Matrix Spike/Matrix Spike Duplicates	A	
VII.	Duplicate sample analysis	A	
VIII.	Laboratory control samples	A	LCS
IX.	Field duplicates	SW	(3,5)
X.	Sample result verification	N	
XI	Overall assessment of data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	MW-23-4	1720817-02	Water	07/28/17
2	MW-23-3	1720817-03	Water	07/28/17
3	MW-23-2	1720817-04	Water	07/28/17
4	MW-23-1	1720817-05	Water	07/28/17
5	DUPE -3-3Q17	1720817-06	Water	07/28/17
6	MW-26-2	1720817-07	Water	07/28/17
7	MW-26-1	1720817-08	Water	07/28/17
8	EB-5-072817	1720817-09	Water	07/28/17
9	MW-23-4MS	1720817-02MS	Water	07/28/17
10	MW-23-4MSD	1720817-02MSD	Water	07/28/17
11	MW-23-4DUP	1720817-02DUP	Water	07/28/17
12				
13				
14				
15				
16				

Notes: _____

VALIDATION FINDINGS WORKSHEET
Sample Specific Analysis Reference

All circled methods are applicable to each sample.

Sample ID	Parameter
1	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
2-8	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
OC	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
9-11	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ CIO ₄

Comments: _____

VALIDATION FINDINGS WORKSHEET
Field DuplicatesInorganics, Method See Cover

Analyte	Concentration		RPD	
	3	5		
Perchlorate (ug/L)	3.9	4.7	19	
Hexavalent Chromium (mg/L)	0.0011	0.00099	11	

V:\FIELD DUPLICATES\Field Duplicates\FD_inorganic\2017\39385A6.wpd

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017
LDC Report Date: September 21, 2017
Parameters: Volatiles
Validation Level: Level III & IV
Laboratory: BC Laboratories, Inc.
Sample Delivery Group (SDG): 17-20934

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
TB-6-073117	1720934-01	Water	07/31/17
MW-4-3	1720934-02	Water	07/31/17
MW-4-2	1720934-03	Water	07/31/17
MW-4-1	1720934-04	Water	07/31/17
MW-21-5	1720934-05	Water	07/31/17
MW-21-4	1720934-06	Water	07/31/17
MW-21-3	1720934-07	Water	07/31/17
MW-21-2**	1720934-08**	Water	07/31/17
MW-21-1	1720934-09	Water	07/31/17
EB-6-073117	1720934-10	Water	07/31/17
SB-2-073117	1720934-11	Water	07/31/17
MW-21-3MS	1720934-07MS	Water	07/31/17
MW-21-3MSD	1720934-07MSD	Water	07/31/17

**Indicates sample underwent Level IV review

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Superfund Organic Methods Data Review (June 2008). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following method:

Volatile Organic Compounds (VOCs) by Environmental Protection Agency (EPA) Method 524.2

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results. Samples appended with a double asterisk on the cover page were subjected to Level IV evaluation, which is comprised of the QC summary forms as well as the raw data, to confirm sample quantitation and identification.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition and cooler temperatures upon receipt met validation criteria.

All technical holding time requirements were met.

II. GC/MS Instrument Performance Check

Instrument performance was checked at 12 hour intervals.

All ion abundance requirements were met.

III. Initial Calibration and Initial Calibration Verification

An initial calibration was performed as required by the method.

For compounds where average relative response factors (RRFs) were utilized, the percent relative standard deviations (%RSD) were less than or equal to 20.0%.

In the case where the laboratory used a calibration curve to evaluate the compounds, all coefficients of determination (r^2) were greater than or equal to 0.990.

The percent differences (%D) of the initial calibration verification (ICV) standard were less than or equal to 30.0% for all compounds.

IV. Continuing Calibration

Continuing calibration was performed at the required frequencies.

The percent differences (%D) were less than or equal to 30.0% for all compounds with the following exceptions:

Date	Compound	%D	Associated Samples	Flag	A or P
08/03/17	Bromomethane Methyl iodide Pentachloroethane	50.5 (≤30) 61.7 (≤30) 35.5 (≤30)	TB-6-073117 MW-4-3 MW-4-2 MW-4-1 MW-21-5 MW-21-4 MW-21-3 MW-21-1 EB-6-073117 SB-2-073117	UJ (all non-detects) UJ (all non-detects) UJ (all non-detects)	P
08/04/17	Methyl iodide Pentachloroethane	51.8 (≤30) 33.9 (≤30)	MW-21-2**	UJ (all non-detects) UJ (all non-detects)	P

V. Laboratory Blanks

Laboratory blanks were analyzed as required by the method. No contaminants were found in the laboratory blanks.

VI. Field Blanks

Sample TB-6-073117 was identified as a trip blank. No contaminants were found.

Sample EB-6-073117 was identified as an equipment blank. No contaminants were found.

Sample SB-2-073117 was identified as a source blank. No contaminants were found.

VII. Surrogates

Surrogates were added to all samples as required by the method. All surrogate recoveries (%R) were within QC limits.

VIII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

IX. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the method. Percent recoveries (%R) were within QC limits.

X. Field Duplicates

No field duplicates were identified in this SDG.

XI. Internal Standards

All internal standard areas and retention times were within QC limits.

XII. Compound Quantitation

All compound quantitations met validation criteria for samples which underwent Level IV validation. Raw data were not reviewed for Level III validation.

XIII. Target Compound Identifications

All target compound identifications met validation criteria for samples which underwent Level IV validation. Raw data were not reviewed for Level III validation.

XIV. System Performance

The system performance was acceptable for samples which underwent Level IV validation. Raw data were not reviewed for Level III validation.

XV. Overall Assessment of Data

The analysis was conducted within all specifications of the method. No results were rejected in this SDG.

Due to continuing calibration %D, data were qualified as estimated in ten samples.

The quality control criteria reviewed, other than those discussed above, were met and are considered acceptable. Sample results that were found to be estimated (J) are usable for limited purposes only. Based upon the data validation all other results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Volatiles - Data Qualification Summary - SDG 17-20934

Sample	Compound	Flag	A or P	Reason
TB-6-073117 MW-4-3 MW-4-2 MW-4-1 MW-21-5 MW-21-4 MW-21-3 MW-21-1 EB-6-073117 SB-2-073117	Bromomethane Methyl iodide Pentachloroethane	UJ (all non-detects) UJ (all non-detects) UJ (all non-detects)	P	Continuing calibration (%D)
MW-21-2**	Methyl iodide Pentachloroethane	UJ (all non-detects) UJ (all non-detects)	P	Continuing calibration (%D)

NASA JPL, 3Q2017
Volatiles - Laboratory Blank Data Qualification Summary - SDG 17-20934

No Sample Data Qualified in this SDG



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

TB-6-073117

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-01 File ID: 03AUG17.D
Sampled: 07/31/17 07:00 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 15:37
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>U5</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

N 192117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

TB-6-073117

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-01 File ID: 03AUG17.D
Sampled: 07/31/17 07:00 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 15:37
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

Handwritten signature/initials



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

TB-6-073117

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-01 File ID: 03AUG17.D
Sampled: 07/31/17 07:00 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 15:37
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U UJ
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U UJ
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.580	106	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.8400	98.4	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.6800	96.8	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	200504	6.57	200979	6.57	
Chlorobenzene-d5 (IS)	76910	9.62	63715	9.61	
1,4-Difluorobenzene (IS)	282629	7.38	300872	7.38	

* Values outside of QC limits

8209-117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

TB-6-073117

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-20934</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1720934-01</u>	File ID:	<u>03AUG17.D</u>		
Sampled:	<u>07/31/17 07:00</u>	Prepared:	<u>08/03/17 07:00</u>	Analyzed:	<u>08/03/17 15:37</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>B[H0282</u>	Sequence:	<u>1713705</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-4-3

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-02 File ID: 03AUG18.D
Sampled: 07/31/17 07:35 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 16:00
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>WJ</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/29/17

Tidewater Inc.
 3761 Attucks Drive
 Powell, OH 43065

 Reported: 8/25/2017 12:56:21PM
 Project: JPL- GW Monitoring Wells
 Project Number: 3Q17
 Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-4-3

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-20934</u>
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>
Matrix:	<u>Water</u>	Laboratory ID:	<u>1720934-02</u>
		File ID:	<u>03AUG18.D</u>
Sampled:	<u>07/31/17 07:35</u>	Prepared:	<u>08/03/17 07:00</u>
		Analyzed:	<u>08/03/17 16:00</u>
Solids:		Preparation:	<u>EPA 5030 Water MS</u>
		Initial/Final:	<u>25 ml / 25 ml</u>
Batch:	<u>BIH0282</u>	Sequence:	<u>1713705</u>
		Calibration:	<u>1707017</u>
		Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.70	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/25/17



Tidewater Inc.
 3761 Attucks Drive
 Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
 Project: JPL- GW Monitoring Wells
 Project Number: 3Q17
 Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
 EPA-524.2

MW-4-3

Laboratory: BC Laboratories SDG: 17-20934
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720934-02 File ID: 03AUG18.D
 Sampled: 07/31/17 07:35 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 16:00
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	11.180	112	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.6500	96.5	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.4600	94.6	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	193902	6.57	200979	6.57	
Chlorobenzene-d5 (IS)	78487	9.62	63715	9.61	
1,4-Difluorobenzene (IS)	286480	7.38	300872	7.38	

* Values outside of QC limits

809247



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-4-3

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-20934</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1720934-02</u>	File ID:	<u>03AUG18.D</u>		
Sampled:	<u>07/31/17 07:35</u>	Prepared:	<u>08/03/17 07:00</u>	Analyzed:	<u>08/03/17 16:00</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0282</u>	Sequence:	<u>1713705</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/25/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-4-2

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-03 File ID: 03AUG19.D
Sampled: 07/31/17 08:05 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 16:23
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>45</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.37	J
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8209217



Tidewater Inc. Reported: 8/25/2017 12:56:21PM
 3761 Attucks Drive Project: JPL- GW Monitoring Wells
 Powell, OH 43065 Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-4-2

Laboratory: BC Laboratories SDG: 17-20934
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720934-03 File ID: 03AUG19.D
 Sampled: 07/31/17 08:05 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 16:23
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U UT
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U US
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	11.310	113	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.7000	97.0	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.6400	96.4	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	191181	6.57	200979	6.57	
Chlorobenzene-d5 (IS)	76438	9.62	63715	9.61	
1,4-Difluorobenzene (IS)	283280	7.38	300872	7.38	

* Values outside of QC limits

209217

Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-4-2

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-03 File ID: 03AUG19.D
Sampled: 07/31/17 08:05 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 16:23
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

2092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-4-1

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-04 File ID: 03AUG20.D
Sampled: 07/31/17 08:30 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 16:46
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>LS</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

2052117



BC Laboratories, Inc.

Environmental Testing Laboratory Since 1949

Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-4-1

Laboratory:	BC Laboratories	SDG:	17-20934
Client:	Tidewater Inc.	Project:	JPL- GW Monitoring Wells
Matrix:	Water	Laboratory ID:	1720934-04
Sampled:	07/31/17 08:30	Prepared:	08/03/17 07:00
Solids:		Preparation:	EPA 5030 Water MS
Batch:	BIH0282	Sequence:	1713705
		Calibration:	1707017
		Instrument:	MS-V5
		File ID:	03AUG20.D
		Analyzed:	08/03/17 16:46
		Initial/Final:	25 ml / 25 ml

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-4-1

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-04 File ID: 03AUG20.D
Sampled: 07/31/17 08:30 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 16:46
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U UJ
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U UJ
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	11.280	113	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.5900	95.9	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.7700	97.7	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	165465	6.57	200979	6.57	
Chlorobenzene-d5 (IS)	68152	9.61	63715	9.61	
1,4-Difluorobenzene (IS)	245587	7.38	300872	7.38	

* Values outside of QC limits

2092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-21-5

Laboratory: BC Laboratories SDG: 17-20934
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720934-05 File ID: 03AUG21.D
 Sampled: 07/31/17 10:15 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 17:09
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.95	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-5

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-05 File ID: 03AUG21.D
Sampled: 07/31/17 10:15 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 17:09
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U UJ
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U UJ
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	11.730	117	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.6700	96.7	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.7900	97.9	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	187987	6.58	200979	6.57	
Chlorobenzene-d5 (IS)	74057	9.61	63715	9.61	
1,4-Difluorobenzene (IS)	280520	7.38	300872	7.38	

* Values outside of QC limits

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-5

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-05 File ID: 03AUG21.D
Sampled: 07/31/17 10:15 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 17:09
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/25/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-4

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-06 File ID: 03AUG22.D
Sampled: 07/31/17 09:45 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 17:32
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U 45
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	5.6	
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-4

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-06 File ID: 03AUG22.D
Sampled: 07/31/17 09:45 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 17:32
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	1.2	
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.30	J
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

259217



Tidewater Inc. Reported: 8/25/2017 12:56:21PM
 3761 Attucks Drive Project: JPL- GW Monitoring Wells
 Powell, OH 43065 Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-4

Laboratory: BC Laboratories SDG: 17-20934
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720934-06 File ID: 03AUG22.D
 Sampled: 07/31/17 09:45 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 17:32
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BJH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U UJ
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U UJ
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.540	105	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.7300	97.3	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	11.510	115	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	189899	6.58	200979	6.57	
Chlorobenzene-d5 (IS)	65647	9.61	63715	9.61	
1,4-Difluorobenzene (IS)	277582	7.38	300872	7.38	

* Values outside of QC limits

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-4

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-06 File ID: 03AUG22.D
Sampled: 07/31/17 09:45 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 17:32
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-3

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-07 File ID: 03AUG08.D
Sampled: 07/31/17 10:50 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 12:10
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>UJ</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.47	J
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.17	J
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

2092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-3

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-07</u>	File ID: <u>03AUG08.D</u>	
Sampled: <u>07/31/17 10:50</u>	Prepared: <u>08/03/17 07:00</u>	Analyzed: <u>08/03/17 12:10</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BH0282</u>	Sequence: <u>1713705</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	1.1	
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	1.1	
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

SLG 8/21/17



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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-3

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-07 File ID: 03AUG08.D
Sampled: 07/31/17 10:50 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 12:10
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.650	106	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.8500	98.5	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	10.060	101	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	192456	6.57	200979	6.57	
Chlorobenzene-d5 (IS)	75222	9.61	63715	9.61	
1,4-Difluorobenzene (IS)	290157	7.38	300872	7.38	

* Values outside of QC limits

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-3

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-20934</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1720934-07</u>	File ID:	<u>03AUG08.D</u>		
Sampled:	<u>07/31/17 10:50</u>	Prepared:	<u>08/03/17 07:00</u>	Analyzed:	<u>08/03/17 12:10</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BJH0282</u>	Sequence:	<u>1713705</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

SC 8/21/17



Laboratories, Inc.

Environmental Testing Laboratory Since 1949

Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-2

Laboratory:	BC Laboratories	SDG:	17-20934
Client:	Tidewater Inc.	Project:	JPL- GW Monitoring Wells
Matrix:	Water	Laboratory ID:	1720934-08
		File ID:	04AUG08.D
Sampled:	07/31/17 11:30	Prepared:	08/03/17 07:00
		Analyzed:	08/04/17 11:59
Solids:		Preparation:	EPA 5030 Water MS
		Initial/Final:	25 ml / 25 ml
Batch:	B/H0282	Sequence:	1713791
		Calibration:	1707017
		Instrument:	MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.16	J
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-08</u>	File ID: <u>04AUG08.D</u>	
Sampled: <u>07/31/17 11:30</u>	Prepared: <u>08/03/17 07:00</u>	Analyzed: <u>08/04/17 11:59</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BIH0282</u>	Sequence: <u>1713791</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.56	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

SL092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-2

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-08 File ID: 04AUG08.D
Sampled: 07/31/17 11:30 Prepared: 08/03/17 07:00 Analyzed: 08/04/17 11:59
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0282 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U UT
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U UT
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	9.1000	91.0	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.9900	99.9	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	10.120	101	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	210639	6.57	230753	6.57	
Chlorobenzene-d5 (IS)	87620	9.62	81486	9.61	
1,4-Difluorobenzene (IS)	300855	7.38	338008	7.38	

* Values outside of QC limits

8/20/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-21-2

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-08 File ID: 04AUG08.D
Sampled: 07/31/17 11:30 Prepared: 08/03/17 07:00 Analyzed: 08/04/17 11:59
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0282 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/29/17



Tidewater Inc. Reported: 8/25/2017 12:56:21PM
 3761 Attucks Drive Project: JPL- GW Monitoring Wells
 Powell, OH 43065 Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-1

Laboratory: BC Laboratories SDG: 17-20934
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720934-09 File ID: 03AUG24.D
 Sampled: 07/31/17 12:00 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 18:18
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: B[H0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U UJ
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.61	
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-09</u>	File ID: <u>03AUG24.D</u>	
Sampled: <u>07/31/17 12:00</u>	Prepared: <u>08/03/17 07:00</u>	Analyzed: <u>08/03/17 18:18</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>B[H0282</u>	Sequence: <u>1713705</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.20	J
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-1

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-09 File ID: 03AUG24.D
Sampled: 07/31/17 12:00 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 18:18
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	11.010	110	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.8900	98.9	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	10.340	103	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	188879	6.58	200979	6.57	
Chlorobenzene-d5 (IS)	71207	9.62	63715	9.61	
1,4-Difluorobenzene (IS)	275631	7.38	300872	7.38	

* Values outside of QC limits

SL57217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-21-1

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-20934</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1720934-09</u>	File ID:	<u>03AUG24.D</u>		
Sampled:	<u>07/31/17 12:00</u>	Prepared:	<u>08/03/17 07:00</u>	Analyzed:	<u>08/03/17 18:18</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BH0282</u>	Sequence:	<u>1713705</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-6-073117

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-10</u>	File ID: <u>03AUG25.D</u>	
Sampled: <u>07/31/17 12:15</u>	Prepared: <u>08/03/17 07:00</u>	Analyzed: <u>08/03/17 18:41</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BJH0282</u>	Sequence: <u>1713705</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>UJ</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/25/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-6-073117

Laboratory: BC Laboratories SDG: 17-20934
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720934-10 File ID: 03AUG25.D
 Sampled: 07/31/17 12:15 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 18:41
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

Scopm17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-6-073117

Laboratory: BC Laboratories SDG: 17-20934
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1720934-10 File ID: 03AUG25.D
Sampled: 07/31/17 12:15 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 18:41
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UT</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UT</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	11.100	111	75 - 125	
Toluene-d8 (Surrogate)	10.000	10.010	100	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.8800	98.8	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	185070	6.57	200979	6.57	
Chlorobenzene-d5 (IS)	75442	9.62	63715	9.61	
1,4-Difluorobenzene (IS)	276520	7.38	300872	7.38	

* Values outside of QC limits

209217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-6-073117

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-10</u>	File ID: <u>03AUG25.D</u>	
Sampled: <u>07/31/17 12:15</u>	Prepared: <u>08/03/17 07:00</u>	Analyzed: <u>08/03/17 18:41</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BH0282</u>	Sequence: <u>1713705</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

SB-2-073117

Laboratory: BC Laboratories SDG: 17-20934
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720934-11 File ID: 03AUG26.D
 Sampled: 07/31/17 12:30 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 19:04
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>U5</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:56:21PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

SB-2-073117

Laboratory: BC Laboratories SDG: 17-20934
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720934-11 File ID: 03AUG26.D
 Sampled: 07/31/17 12:30 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 19:04
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

SL 082117



Tidewater Inc. Reported: 8/25/2017 12:56:21PM
 3761 Attucks Drive Project: JPL- GW Monitoring Wells
 Powell, OH 43065 Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

SB-2-073117

Laboratory: BC Laboratories SDG: 17-20934
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1720934-11 File ID: 03AUG26.D
 Sampled: 07/31/17 12:30 Prepared: 08/03/17 07:00 Analyzed: 08/03/17 19:04
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0282 Sequence: 1713705 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	11.070	111	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.9100	99.1	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.9000	99.0	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	189426	6.57	200979	6.57	
Chlorobenzene-d5 (IS)	75627	9.61	63715	9.61	
1,4-Difluorobenzene (IS)	275283	7.38	300872	7.38	

* Values outside of QC limits

8/20/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:56:21PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

SB-2-073117

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-20934</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1720934-11</u>	File ID:	<u>03AUG26.D</u>		
Sampled:	<u>07/31/17 12:30</u>	Prepared:	<u>08/03/17 07:00</u>	Analyzed:	<u>08/03/17 19:04</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0282</u>	Sequence:	<u>1713705</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

209217

LDC #: 39385B1

VALIDATION COMPLETENESS WORKSHEET

SDG #: 17-20934

Level III/IV

Laboratory: BC Laboratories, Inc.

Date: 9/12/17

Page: 1 of 1

Reviewer: [Signature]

2nd Reviewer: [Signature]

METHOD: GC/MS Volatiles (EPA Method 524.2)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A	
II.	GC/MS Instrument performance check	A	
III.	Initial calibration/ICV	AA	RSD ≤ 20%. Y ² ICV ≤ 30%
IV.	Continuing calibration	MW	CCV ≤ 30%
V.	Laboratory Blanks	A	
VI.	Field blanks	ND	TB=1. EB=10. SB=11
VII.	Surrogate spikes	A	
VIII.	Matrix spike/Matrix spike duplicates	A	
IX.	Laboratory control samples	A	ICS
X.	Field duplicates	N	
XI.	Internal standards	A	
XII.	Compound quantitation RL/LOQ/LODs	A	Not reviewed for Level III validation.
XIII.	Target compound identification	A	Not reviewed for Level III validation.
XIV.	System performance	A	Not reviewed for Level III validation.
XV.	Overall assessment of data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

** Indicates sample was underwent Level IV review

	Client ID	Lab ID	Matrix	Date
1	TB-6-073117	1720934-01	Water	07/31/17
2	MW-4-3	1720934-02	Water	07/31/17
3	MW-4-2	1720934-03	Water	07/31/17
4	MW-4-1	1720934-04	Water	07/31/17
5	MW-21-5	1720934-05	Water	07/31/17
6	MW-21-4	1720934-06	Water	07/31/17
7	MW-21-3	1720934-07	Water	07/31/17
8	MW-21-2**	1720934-08**	Water	07/31/17
9	MW-21-1	1720934-09	Water	07/31/17
10	EB-6-073117	1720934-10	Water	07/31/17
11	SB-2-073117	1720934-11	Water	07/31/17
12	MW-21-3MS	1720934-07MS	Water	07/31/17
13	MW-21-3MSD	1720934-07MSD	Water	07/31/17

Method: Volatiles (EPA Method 524.2)

Validation Area	Yes	No	NA	Findings/Comments
I. Technical holding times				
Were all technical holding times met?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was cooler temperature criteria met?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
II. GC/MS Instrument performance check				
Was a tune check performed prior to establishing and/or re-establishing an initial calibration?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the BFB performance results reviewed and found to be within the specified criteria?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
III. Initial calibration				
Did the laboratory perform at least 5 point calibration prior to sample analysis?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all percent relative standard deviations (%RSD) \leq 20%?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
IIIa. Initial Calibration Verification calibration				
Was an initial calibration verification standard analyzed after each initial calibration for each instrument?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all percent differences (%D) \leq 30%?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
IV. Continuing calibration				
Was a continuing calibration standard analyzed at the beginning of each analysis batch?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all percent differences (%D) of continuing calibration $<$ 30%?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
V. Laboratory Blanks				
Was a laboratory blank associated with every sample in this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was a laboratory blank analyzed with each analysis batch?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was there contamination in the laboratory blanks? If yes, please see the Blanks validation completeness worksheet.	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
VI. Field blanks				
Field blanks were identified in this SDG.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Target compounds were detected in the field blanks.	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
VII. Surrogate spikes				
Were all surrogate %R within the QC limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
If the percent recovery (%R) for one or more surrogates was out of QC limits, was a reanalysis performed to confirm samples with %R outside of criteria?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
VIII. Matrix spike/Matrix spike duplicates				
Was a matrix spike (MS) and matrix spike duplicate (MSD) analyzed for this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the QC limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
IX. Laboratory control samples				
Was an LCS analyzed for this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	

Validation Area	Yes	No	NA	Findings/Comments
Was an LCS analyzed per analytical batch?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the LCS percent recoveries (%R) within 70-130%?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
X. Field duplicates				
Field duplicate pairs were identified in this SDG.	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
Target compounds were detected in the field duplicates.	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
XI. Internal standards				
Were internal standard area counts within +/-30% of the area of the most recent continuing calibration standard and +/-50% of the average peak area in the initial calibration?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were retention times within +/-30 seconds of the associated calibration standard?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
XII. Compound quantitation/CRQLs				
Were the correct internal standard (IS), quantitation ion and relative response factor (RRF) or regression equations used to quantitate the compound?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were compound quantitation and CRQLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
XIII. Target compound identification				
Were relative retention times (RRT's) within + 0.06 RRT units of the standard?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Did compound spectra meet specified EPA "Functional Guidelines" criteria?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were chromatogram peaks verified and accounted for?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
XIV. System performance				
System performance was found to be acceptable.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
XV. Overall assessment of data				
Overall assessment of data was found to be acceptable.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	

TARGET COMPOUND WORKSHEET

METHOD: VOA

A. Chloromethane	AA. Tetrachloroethene	AAA. 1,3,5-Trimethylbenzene	AAAA. Ethyl tert-butyl ether	A1. 1,3-Butadiene
B. Bromomethane	BB. 1,1,2,2-Tetrachloroethane	BBB. 4-Chlorotoluene	BBBB. tert-Amyl methyl ether	B1. Hexane
C. Vinyl chloride	CC. Toluene	CCC. tert-Butylbenzene	CCCC. 1-Chlorohexane	C1. Heptane
D. Chloroethane	DD. Chlorobenzene	DDD. 1,2,4-Trimethylbenzene	DDDD. Isopropyl alcohol	D1. Propylene
E. Methylene chloride	EE. Ethylbenzene	EEE. sec-Butylbenzene	EEEE. Acetonitrile	E1. Freon 11
F. Acetone	FF. Styrene	FFF. 1,3-Dichlorobenzene	FFFF. Acrolein	F1. Freon 12
G. Carbon disulfide	GG. Xylenes, total	GGG. p-Isopropyltoluene	GGGG. Acrylonitrile	G1. Freon 113
H. 1,1-Dichloroethene	HH. Vinyl acetate	HHH. 1,4-Dichlorobenzene	HHHH. 1,4-Dioxane	H1. Freon 114
I. 1,1-Dichloroethane	II. 2-Chloroethylvinyl ether	III. n-Butylbenzene	IIII. Isobutyl alcohol	I1. 2-Nitropropane
J. 1,2-Dichloroethene, total	JJ. Dichlorodifluoromethane	JJJ. 1,2-Dichlorobenzene	JJJJ. Methacrylonitrile	J1. Dimethyl disulfide
K. Chloroform	KK. Trichlorofluoromethane	KKK. 1,2,4-Trichlorobenzene	KKKK. Propionitrile	K1. 2,3-Dimethyl pentane
L. 1,2-Dichloroethane	LL. Methyl-tert-butyl ether	LLL. Hexachlorobutadiene	LLLL. Ethyl ether	L1. 2,4-Dimethyl pentane
M. 2-Butanone	MM. 1,2-Dibromo-3-chloropropane	MMM. Naphthalene	MMMM. Benzyl chloride	M1. 3,3-Dimethyl pentane
N. 1,1,1-Trichloroethane	NN. Methyl ethyl ketone	NNN. 1,2,3-Trichlorobenzene	NNNN. Iodomethane	N1. 2-Methylpentane
O. Carbon tetrachloride	OO. 2,2-Dichloropropane	OOO. 1,3,5-Trichlorobenzene	OOOO. 1,1-Difluoroethane	O1. 3-Methylpentane
P. Bromodichloromethane	PP. Bromochloromethane	PPP. trans-1,2-Dichloroethene	PPPP. Tetrahydrofuran	P1. 3-Ethylpentane
Q. 1,2-Dichloropropane	QQ. 1,1-Dichloropropene	QQQ. cis-1,2-Dichloroethene	QQQQ. Methyl acetate	Q1. 2,2-Dimethylpentane
R. cis-1,3-Dichloropropene	RR. Dibromomethane	RRR. m,p-Xylenes	RRRR. Ethyl acetate	R1. 2,2,3-Trimethylbutane
S. Trichloroethene	SS. 1,3-Dichloropropane	SSS. o-Xylene	SSSS. Cyclohexane	S1. 2,2,4-Trimethylpentane
T. Dibromochloromethane	TT. 1,2-Dibromoethane	TTT. 1,1,2-Trichloro-1,2,2-trifluoroethane	TTTT. Methylcyclohexane	T1. 2-Methylhexane
U. 1,1,2-Trichloroethane	UU. 1,1,1,2-Tetrachloroethane	UUU. 1,2-Dichlorotetrafluoroethane	UUUU. Allyl chloride	U1. Nonanal
V. Benzene	VV. Isopropylbenzene	VVV. 4-Ethyltoluene	VVVV. Methyl methacrylate	V1. 2-Methylnaphthalene
W. trans-1,3-Dichloropropene	WW. Bromobenzene	WWW. Ethanol	WWWWW. Ethyl methacrylate	W1. Methanol
X. Bromoform	XX. 1,2,3-Trichloropropane	XXX. Di-isopropyl ether	XXXX. cis-1,4-Dichloro-2-butene	X1. 1,2,3-Trimethylbenzene
Y. 4-Methyl-2-pentanone	YY. n-Propylbenzene	YYY. tert-Butanol	YYYY. trans-1,4-Dichloro-2-butene	Y1.
Z. 2-Hexanone	ZZ. 2-Chlorotoluene	ZZZ. tert-Butyl alcohol	ZZZZ. Pentachloroethane	Z1.

VALIDATION FINDINGS WORKSHEET

Continuing Calibration

METHOD: GC/MS VOA (EPA Method 524.2)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

N N/A Was a continuing calibration standard analyzed at least once every 12 hours for each instrument?

Y N N/A Were all percent differences (%D) < 30% ?

#	Date	Standard ID	Compound	Finding %D (Limit: <30.0%)	Associated Samples	Qualifications
	<u>8/3/17</u>	<u>03AUG02</u>	<u>B</u> <u>Methyl iodide</u>	<u>50.5</u> <u>61.7</u>	<u>1-7. 9-13. MB</u> <u>(ND)</u>	<u>✓/N/A</u> <u>↓</u>
			<u>2222</u>	<u>35.5</u>		
	<u>8/4/17</u>	<u>04AUG03</u>	<u>Methyl iodide</u>	<u>51.8</u> <u>33.9</u>	<u>8. MB (ND)</u>	<u>✓/N/A</u> <u>↓</u>
			<u>2222</u>			

VALIDATION FINDINGS WORKSHEET Initial Calibration Calculation Verification

METHOD: GC/MS VOA (EPA SW 846 Method 8260B)

The Relative Response Factor (RRF), average RRF, and percent relative standard deviation (%RSD) were recalculated for the compounds identified below using the following calculations:

$$RRF = (A_x)(C_{is}) / (A_{is})(C_x)$$

average RRF = sum of the RRFs/number of standards

$$\%RSD = 100 * (S/X)$$

A_x = Area of compound,

C_x = Concentration of compound,

S = Standard deviation of the RRFs

X = Mean of the RRFs

A_{is} = Area of associated internal standard

C_{is} = Concentration of internal standard

#	Standard ID	Calibration Date	Compound (Reference Internal Standard)	Reported	Recalculated	Reported	Recalculated	Reported	Recalculated
				RRF (10 std)	RRF (10 std)	Average RRF (initial)	Average RRF (initial)	%RSD	%RSD
1	ICAL (V5)	7/20/17	K (1st internal standard)	0.7797252	0.7797252	0.7571101	0.7571101	5.641054	5.641053
			S (2nd internal standard)	0.3603684	0.3603684	0.3434012	0.3434012	3.079274	3.079306
			EE (3rd internal standard)	2.063766	2.063766	1.948304	1.948304	6.711166	6.711173
			BB (4th internal standard)						
2			K (1st internal standard)						
			S (2nd internal standard)						
			AA (3rd internal standard)						
			BB (4th internal standard)						
3			K (1st internal standard)						
			S (2nd internal standard)						
			AA (3rd internal standard)						
			BB (4th internal standard)						
4			K (1st internal standard)						
			S (2nd internal standard)						
			AA (3rd internal standard)						
			BB (4th internal standard)						

Comments: Refer to Initial Calibration findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

VALIDATION FINDINGS WORKSHEET Continuing Calibration Results Verification

METHOD: GC/MS VOA (EPA Method 524.2)

The percent difference (%D) of the initial calibration average Relative Response Factors (RRFs) and the continuing calibration RRFs were recalculated for the compounds identified below using the following calculation:

% Difference = 100 * (ave. RRF - RRF)/ave. RRF
 RRF = (A_x)(C_{is})/(A_{is})(C_x)

Where: ave. RRF = initial calibration average RRF
 RRF = continuing calibration RRF
 A_x = Area of compound, A_{is} = Area of associated internal standard
 C_x = Concentration of compound, C_{is} = Concentration of internal standard

#	Standard ID	Calibration Date	Compound (Reference internal Standard)	Average RRF (initial)	Reported RRF (CC)	Recalculated RRF (CC)	Reported %D	Recalculated %D
1	04aug02	8/4/17	K (1st internal standard)	0.7571101	0.7733756	0.7733756	2.1	2.1
			S (2nd internal standard)	0.3434012	0.3649989	0.3649989	6.3	6.3
			EE (3rd internal standard)	1.948304	2.011687	2.011687	3.3	3.3
			BB (4th internal standard)					
2			K (5th internal standard)					
			S (2nd internal standard)					
			AA (3rd internal standard)					
			BB (4th internal standard)					
3			QQQ (1st internal standard)					
			S (2nd internal standard)					
			AA (3rd internal standard)					
			MMM (4th internal standard)					
4			BB (5th internal standard)					
			S (2nd internal standard)					
			AA (3rd internal standard)					
			BB (4th internal standard)					

Comments: Refer to Continuing Calibration findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 29325B1

VALIDATION FINDINGS WORKSHEET
Surrogate Results Verification

Page: 1 of 1
Reviewer: [Signature]
2nd reviewer: [Signature]

METHOD: GC/MS VOA (EPA Method 524.2)

The percent recoveries (%R) of surrogates were recalculated for the compounds identified below using the following calculation:

% Recovery: SF/SS * 100

Where: SF = Surrogate Found
SS = Surrogate Spiked

Sample ID: 8

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8	10.00	9.99	99.9	99.9	0
Bromofluorobenzene	↓	10.12	101	101	↓
1,2-Dichlorobenzene-d4	↓	9.10	91.0	91.0	↓
Dibromofluoromethane					

Sample ID:

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					
Dibromofluoromethane					

Sample ID:

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					
Dibromofluoromethane					

Sample ID:

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					
Dibromofluoromethane					

LDC #: 29385B

VALIDATION FINDINGS WORKSHEET

Matrix Spike/Matrix Spike Duplicates Results Verification

Page: 1 of 1
 Reviewer: Q
 2nd Reviewer: KW

METHOD: GC/MS VOA (EPA Method 524.2)

The percent recoveries (%R) and Relative Percent Difference (RPD) of the matrix spike and matrix spike duplicate were recalculated for the compounds identified below using the following calculation:

% Recovery = $100 * (SSC - SC) / SA$

Where: SSC = Spiked sample concentration
 SA = Spike added

SC = Sample concentration

RPD = $|MSC - MSC| * 2 / (MSC + MSDC)$

MSC = Matrix spike concentration

MSDC = Matrix spike duplicate concentration

MS/MSD sample: 12/13

Compound	Spike Added		Sample Concentration	Spiked Sample Concentration		Matrix Spike		Matrix Spike Duplicate		MS/MSD	
	MS	MSD		MS	MSD	Percent Recovery		Percent Recovery		RPD	
						Reported	Recalc	Reported	Recalc	Reported	Recalculated
1,1-Dichloroethene	25.00	25.00	ND	30.780	27.270	123	123	109	109	12.1	12.1
Trichloroethene	↓	↓	1.06	30.140	27.520	116	116	106	106	9.09	9.09
Benzene	↓	↓	ND	25.920	23.420	104	104	93.7	93.7	10.1	10.1
Toluene	↓	↓	↓	27.270	26.260	109	109	105	105	3.77	3.77
Chlorobenzene	↓	↓	↓	24.370	23.200	97.5	97.5	92.8	92.8	4.92	4.92

Comments: Refer to Matrix Spike/Matrix Spike Duplicates findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

VALIDATION FINDINGS WORKSHEET
Laboratory Control Sample Results Verification

METHOD: GC/MS VOA (EPA Method 524.2)

The percent recoveries (%R) and Relative Percent Difference (RPD) of the laboratory control sample and laboratory control sample duplicate (if applicable) were recalculated for the compounds identified below using the following calculation:

% Recovery = 100 * SSC/SA

Where: SSC = Spiked sample concentration
 SA = Spike added

RPD = | LCSC - LCSDC | * 2 / (LCSC + LCSDC)

LCSC = Laboratory control sample concentration LCSDC = Laboratory control sample duplicate concentration

LCS ID: BH 0202-B51

Compound	Spike Added (µg/L)		Spiked Sample Concentration (µg/L)		LCS		LCSD		LCS/LCSD	
	LCS	LCSD	LCS	LCSD	Percent Recovery		Percent Recovery		RPD	
					Reported	Recalc.	Reported	Recalc.	Reported	Recalculated
1,1-Dichloroethene	25.000	NA	28.970	NA	116	116				
Trichloroethene	↓	↓	27.730	↓	111	111				
Benzene	↓	↓	24.330	↓	97.3	97.3				
Toluene	↓	↓	26.730	↓	107	107				
Chlorobenzene	↓	↓	25.180	↓	101	101				

Comments: Refer to Laboratory Control Sample findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

VALIDATION FINDINGS WORKSHEET Sample Calculation Verification

METHOD: GC/MS VOA (EPA Method 524.2)

Y N N/A
Y N N/A

Were all reported results recalculated and verified for all level IV samples?

Were all recalculated results for detected target compounds agree within 10.0% of the reported results?

$$\text{Concentration} = \frac{(A_x)(I_s)(DF)}{(A_{is})(RRF)(V_o)(\%S)}$$

- A_x = Area of the characteristic ion (EICP) for the compound to be measured
- A_{is} = Area of the characteristic ion (EICP) for the specific internal standard
- I_s = Amount of internal standard added in nanograms (ng)
- RRF = Relative response factor of the calibration standard.
- V_o = Volume or weight of sample pruged in milliliters (ml) or grams (g).
- Df = Dilution factor.
- %S = Percent solids, applicable to soils and solid matrices only.

Example:

Sample I.D. 8, ~~AA~~

$$\text{Conc} = \frac{1015 (10.0) (1)}{30085 (0.329699) (1)}$$

$$\text{Conc} = \frac{(5520) (10.0) (1)}{(87620) (30085) (0.329699)} = 0.56 \mu\text{g/L}$$

#	Sample ID	Compound	Reported Concentration ($\mu\text{g/L}$)	Calculated Concentration ()	Qualification
	8	AA	0.56		

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017

LDC Report Date: September 20, 2017

Parameters: Chromium

Validation Level: Level III

Laboratory: BC Laboratories, Inc.

Sample Delivery Group (SDG): 17-20934

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-4-3	1720934-02	Water	07/31/17
MW-4-2	1720934-03	Water	07/31/17
MW-4-1	1720934-04	Water	07/31/17
MW-21-5	1720934-05	Water	07/31/17
MW-21-4	1720934-06	Water	07/31/17
MW-21-3	1720934-07	Water	07/31/17
MW-21-2**	1720934-08**	Water	07/31/17
MW-21-1	1720934-09	Water	07/31/17
EB-6-073117	1720934-10	Water	07/31/17
SB-2-073117	1720934-11	Water	07/31/17
MW-21-3MS	1720934-07MS	Water	07/31/17
MW-21-3MSD	1720934-07MSD	Water	07/31/17
MW-21-3DUP	1720934-07DUP	Water	07/31/17

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following method:

Chromium by Environmental Protection Agency (EPA) Method 200.8

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. ICPMS Tune

The mass calibration was within 0.1 AMU and the percent relative standard deviation (%RSD) was less than or equal to 5%.

III. Instrument Calibration

Initial and continuing calibrations were performed as required by the method.

The initial calibration verification (ICV) and continuing calibration verification (CCV) standards were within QC limits.

IV. ICP Interference Check Sample Analysis

ICP interference check sample analyses were not required by the method.

V. Laboratory Blanks

Laboratory blanks were analyzed as required by the method. No contaminants were found in the laboratory blanks with the following exceptions:

Blank ID	Analyte	Maximum Concentration	Associated Samples
PB (prep blank)	Chromium	0.72300 ug/L	All samples in SDG 17-20934

Data qualification by the laboratory blanks was based on the maximum contaminant concentration in the laboratory blanks in the analysis of each analyte. The sample concentrations were either not detected or were significantly greater (>5X blank contaminants) than the concentrations found in the associated laboratory blanks with the following exceptions:

Sample	Analyte	Reported Concentration	Modified Final Concentration
MW-4-1	Chromium	1.8 ug/L	1.8U ug/L
MW-21-5	Chromium	2.6 ug/L	2.6U ug/L
MW-21-4	Chromium	3.0 ug/L	3.0U ug/L
MW-21-2**	Chromium	1.8 ug/L	1.8U ug/L

Sample	Analyte	Reported Concentration	Modified Final Concentration
MW-21-1	Chromium	3.1 ug/L	3.1U ug/L
EB-6-073117	Chromium	2.3 ug/L	2.3U ug/L
SB-2-073117	Chromium	1.9 ug/L	1.9U ug/L

VI. Field Blanks

Sample EB-6-073117 was identified as an equipment blank. No contaminants were found with the following exceptions:

Blank ID	Analyte	Concentration
EB-6-073117	Chromium	2.3 ug/L

Sample SB-2-073117 was identified as an equipment blank. No contaminants were found with the following exceptions:

Blank ID	Analyte	Concentration
SB-2-073117	Chromium	1.9 ug/L

VII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VIII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

IX. Serial Dilution

Serial dilution was not performed for this SDG.

X. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the method. Percent recoveries (%R) were within QC limits.

XI. Field Duplicates

No field duplicates were identified in this SDG.

XII. Internal Standards (ICP-MS)

Raw data were not reviewed for Level III validation.

XIII. Sample Result Verification

Raw data were not reviewed for Level III validation.

XIV. Overall Assessment of Data

The analysis was conducted within all specifications of the method. No results were rejected in this SDG.

Due to laboratory blank contamination, data were qualified as not detected in seven samples.

The quality control criteria reviewed, other than those discussed above, were met and are considered acceptable. Based upon the data validation all other results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Chromium - Data Qualification Summary - SDG 17-20934

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017
Chromium - Laboratory Blank Data Qualification Summary - SDG 17-20934

Sample	Analyte	Modified Final Concentration	A or P
MW-4-1	Chromium	1.8U ug/L	A
MW-21-5	Chromium	2.6U ug/L	A
MW-21-4	Chromium	3.0U ug/L	A
MW-21-2**	Chromium	1.8U ug/L	A
MW-21-1	Chromium	3.1U ug/L	A
EB-6-073117	Chromium	2.3U ug/L	A
SB-2-073117	Chromium	1.9U ug/L	A



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 1:02:11PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
--	---

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-4-3

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-02

File ID: PE-EL3 170803R-168

Sampled: 07/31/17 07:35

Prepared: 08/03/17 08:00

Analyzed: 08/04/17 02:28

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0275

Sequence: 1713865

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	8.6	1		EPA-200.8

11/9/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 1:02:11PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-4-2

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-03

File ID: PE-EL3 170803R-169

Sampled: 07/31/17 08:05

Prepared: 08/03/17 08:00

Analyzed: 08/04/17 02:31

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0275

Sequence: 1713865

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	4.2	1		EPA-200.8

Handwritten: R092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 1:02:11PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-4-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-04</u>	File ID: <u>PE-EL3_170803R-170</u>	
Sampled: <u>07/31/17 08:30</u>	Prepared: <u>08/03/17 08:00</u>	Analyzed: <u>08/04/17 02:34</u>	
Solids: <u>0.00</u>	Preparation: <u>EPA 200.2</u>	Initial/Final: <u>50 ml / 50 ml</u>	
Batch: <u>B[H0275]</u>	Sequence: <u>1713865</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>PE-EL3</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.8 <i>U</i>	1	J	EPA-200.8

8092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 1:02:11PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-21-5

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-05</u>
Sampled: <u>07/31/17 10:15</u>	Prepared: <u>08/03/17 08:00</u>
Solids: <u>0.00</u>	Preparation: <u>EPA 200.2</u>
Batch: <u>BJH0275</u>	Sequence: <u>1713865</u>
	Calibration: <u>UNASSIGNED</u>
	Instrument: <u>PE-EL3</u>
	File ID: <u>PE-EL3 170803R-171</u>
	Analyzed: <u>08/04/17 02:38</u>
	Initial/Final: <u>50 ml / 50 ml</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.6 U	1	J	EPA-200.8

520917



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 1:02:11PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-21-4

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-06

File ID: PE-EL3_170803R-172

Sampled: 07/31/17 09:45

Prepared: 08/03/17 08:00

Analyzed: 08/04/17 02:41

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0275

Sequence: 1713865

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	3.0 U	1		EPA-200.8

809117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 1:02:11PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-21-3

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-07

File ID: PE-EL3 170803R-160

Sampled: 07/31/17 10:50

Prepared: 08/03/17 08:00

Analyzed: 08/04/17 02:00

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0275

Sequence: 1713865

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	4.2	1		EPA-200.8

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 1:02:11PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-21-2

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-08

File ID: PE-EL3 170803R-173

Sampled: 07/31/17 11:30

Prepared: 08/03/17 08:00

Analyzed: 08/04/17 02:45

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0275

Sequence: 1713865

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.8 U	1	J	EPA-200.8

8/25/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 1:02:11PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-21-1

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-09

File ID: PE-EL3 170803R-174

Sampled: 07/31/17 12:00

Prepared: 08/03/17 08:00

Analyzed: 08/04/17 02:48

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0275

Sequence: 1713865

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	3.1 U	1		EPA-200.8

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 1:02:11PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

EB-6-073117

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-10

File ID: PE-EL3 170803R-175

Sampled: 07/31/17 12:15

Prepared: 08/03/17 08:00

Analyzed: 08/04/17 02:52

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0275

Sequence: 1713865

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.3 <i>u</i>	1	J	EPA-200.8

Slup 8/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 1:02:11PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

SB-2-073117

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-11

File ID: PE-EL3 170803R-176

Sampled: 07/31/17 12:30

Prepared: 08/03/17 08:00

Analyzed: 08/04/17 02:55

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0275

Sequence:

1713865

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.9 U	1	J	EPA-200.8

Handwritten signature or initials

LDC #: 39385B4a

VALIDATION COMPLETENESS WORKSHEET

Date: 9/18/17

SDG #: 17-20934

Level III

Page: 1 of 1

Laboratory: BC Laboratories, Inc.

Reviewer: B2nd Reviewer: PK

METHOD: Chromium (EPA Method 200.8)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A / A	
II.	ICP/MS Tune	A	
III.	Instrument Calibration	A	
IV.	ICP Interference Check Sample (ICS) Analysis	N	Not Required
V.	Laboratory Blanks	SW	
VI.	Field Blanks	SW	EB=9, SB=10
VII.	Matrix Spike/Matrix Spike Duplicates	A	
VIII.	Duplicate sample analysis	A	< 5x RL; OK by Difference
IX.	Serial Dilution	N	Not performed
X.	Laboratory control samples	A	LCS
XI.	Field Duplicates	N	
XII.	Internal Standard (ICP-MS)	A	
XIII.	Sample Result Verification	N	
XIV.	Overall Assessment of Data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	MW-4-3	1720934-02	Water	07/31/17
2	MW-4-2	1720934-03	Water	07/31/17
3	MW-4-1	1720934-04	Water	07/31/17
4	MW-21-5	1720934-05	Water	07/31/17
5	MW-21-4	1720934-06	Water	07/31/17
6	MW-21-3	1720934-07	Water	07/31/17
7	MW-21-2**	1720934-08**	Water	07/31/17
8	MW-21-1	1720934-09	Water	07/31/17
9	EB-6-073117	1720934-10	Water	07/31/17
10	SB-2-073117	1720934-11	Water	07/31/17
11	MW-21-3MS	1720934-07MS	Water	07/31/17
12	MW-21-3MSD	1720934-07MSD	Water	07/31/17
13	MW-21-3DUP	1720934-07DUP	Water	07/31/17
14				

Notes:

Method:Metals (EPA SW 846 Method 6010/6020/7000)

Validation Area	Yes	No	NA	Findings/Comments
<i>I. Technical holding times</i>				
All technical holding times were met.	✓			
Cooler temperature criteria was met.	✓			
<i>II. ICP/MS Tune</i>				
Were all isotopes in the tuning solution mass resolution within 0.1 amu?	✓			
Were %RSD of isotopes in the tuning solution $\leq 5\%$?	✓			
<i>III. Calibration</i>				
Were all instruments calibrated daily, each set-up time?	✓			
Were the proper number of standards used?	✓			
Were all initial and continuing calibration verification %Rs within the 90-110% (80-120% for mercury) QC limits?	✓			
Were the low standard checks within 70-130%			✓	
Were all initial calibration correlation coefficients within limits as specified by the method?	✓			
<i>IV. Blanks</i>				
Was a method blank associated with every sample in this SDG?	✓			
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.	✓			
<i>V. ICP Interference Check Sample</i>				
Were ICP interference check samples performed daily?			✓	
Were the AB solution percent recoveries (%R) with the 80-120% QC limits?			✓	
<i>VI. Matrix spike/Matrix spike duplicates</i>				
Were a matrix spike (MS) and duplicate (DUP) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD or MS/DUP. Soil / Water.	✓			
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the 75-125 QC limits? If the sample concentration exceeded the spike concentration by a factor of 4 or more, no action was taken.	✓			
Were the MS/MSD or duplicate relative percent differences (RPD) $\leq 20\%$ for waters and $\leq 35\%$ for soil samples? A control limit of $\pm RL$ ($\pm 2X RL$ for soil) was used for samples that were $\leq 5X$ the RL, including when only one of the duplicate sample values were $\leq 5X$ the RL.	✓			
<i>VII. Laboratory control samples</i>				
Was an LCS analyzed for this SDG?	✓			
Was an LCS analyzed per extraction batch?	✓			
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the 80-120% QC limits for water samples and laboratory established QC limits for soils?	✓			

Validation Area	Yes	No	NA	Findings/Comments
VIII. Internal Standards (EPA SW 846 Method 6020/EPA 200.8)				
Were all the percent recoveries (%R) within the 30-120% (6020)/60-125% (200.8) of the intensity of the internal standard in the associated initial calibration?	✓			
If the %Rs were outside the criteria, was a reanalysis performed?			✓	
IX. ICP Serial Dilution				
Was an ICP serial dilution analyzed if analyte concentrations were > 50X the MDL (ICP)/>100X the MDL(ICP/MS)?			✓	
Were all percent differences (%Ds) < 10%?			✓	
Was there evidence of negative interference? If yes, professional judgement will be used to qualify the data.		✓		
X. Sample Result Verification				
Were RLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	✓			
XI. Overall assessment of data				
Overall assessment of data was found to be acceptable.	✓			
XII. Field duplicates				
Field duplicate pairs were identified in this SDG.		✓		
Target analytes were detected in the field duplicates.			✓	
XIII. Field blanks				
Field blanks were identified in this SDG.	✓			
Target analytes were detected in the field blanks.	✓			

VALIDATION FINDINGS WORKSHEET
PB/ICB/CCB QUALIFIED SAMPLES

METHOD: Trace metals (EPA SW 864 Method 6010B/6020/7000)
 Sample Concentration units, unless otherwise noted: ug/L

Soil preparation factor applied: NA
 Associated Samples: All

Analyte	Maximum PB ^a (mg/Kg)	Maximum PB ^a (ug/l)	Maximum ICB/CCB ^a (mg/l)	Action Level	3	4	5	7	8	9	10		
Cr		0.72300		3.615	1.8	2.6	3.0	1.8	3.1	2.3	1.9		

Samples with analyte concentrations within five times the associated ICB, CCB or PB concentration are listed above with the identifications from the Validation Completeness Worksheet. These sample results were qualified as not detected, "U".
 Note : a - The listed analyte concentration is the highest ICB, CCB, or PB detected in the analysis of each element.

LDC #: 39385B4a
SDG #: 17-2034

VALIDATION FINDINGS WORKSHEET Field Blanks

Page: 1 of 1
Reviewer: YB
2nd reviewer: KK

METHOD: Trace Metals (EPA CLP SOW ILM02.1)

Y N N/A Were field blanks identified in this SDG?
Y N N/A Were target analytes detected in the field blanks?

Sample: 9 Field Blank / Trip Blank / Rinsate / Other EB (circle one)

Analyte	Concentration Units (µg/L)
Cr	2.3

Sample: 10 Field Blank / Trip Blank / Rinsate / Other SB (circle one)

Analyte	Concentration Units (µg/L)
Cr	1.9

VALIDATION FINDINGS WORKSHEET
Initial and Continuing Calibration Calculation Verification

METHOD: Trace metals (EPA SW 846 Method 6010/6020/7000)

An initial and continuing calibration verification percent recovery (%R) was recalculated for each type of analysis using the following formula:

$\%R = \frac{\text{Found}}{\text{True}} \times 100$

Where, Found = concentration (in ug/L) of each analyte measured in the analysis of the ICV or CCV solution
 True = concentration (in ug/L) of each analyte in the ICV or CCV source

Standard ID	Type of Analysis	Element	Found (ug/L)	True (ug/L)	Recalculated	Reported	Acceptable (Y/N)
					%R	%R	
	ICP (Initial calibration)						
ICV	ICP/MS (Initial calibration)	Cr	50.062 ug/L	50.000 ug/L	100%	100%	Y
	CVAA (Initial calibration)						
	ICP (Continuing calibration)						
CCV #	ICP/MS (Continuing calibration) 02:21	Cr	39.480 ug/L	40.000 ug/L	99.2%	99.2%	Y
	CVAA (Continuing calibration)						

Comments:

LDC #: 39385B4a
 SDG #: 17-20934

VALIDATION FINDINGS WORKSHEET
Level IV Recalculation Worksheet

Page: 1 of 1
 Reviewer: JB
 2nd Reviewer: YV

METHOD: Trace metals (EPA CLP SOW ILM02.1)

Percent recoveries (%R) for an ICP interference check sample, a laboratory control sample and a matrix spike sample were recalculated using the following formula:

$$\%R = \frac{\text{Found}}{\text{True}} \times 100$$

Where, Found = Concentration of each analyte measured in the analysis of the sample. For the matrix spike calculation, Found = SSR (spiked sample result) - SR (sample result).
 True = Concentration of each analyte in the source.

A sample and duplicate relative percent difference (RPD) was recalculated using the following formula:

$$\text{RPD} = \frac{|S-D|}{(S+D)/2} \times 100$$

Where, S = Original sample concentration
 D = Duplicate sample concentration

An ICP serial dilution percent difference (%D) was recalculated using the following formula:

$$\%D = \frac{|I-SDR|}{I} \times 100$$

Where, I = Initial Sample Result (ug/L)
 SDR = Serial Dilution Result (ug/L) (Instrument Reading x 5)

Sample ID	Type of Analysis	Element	Found / S / I (units)	True / D / SDR (units)	Recalculated	Reported	Acceptable (Y/N)
					%R / RPD / %D	%R / RPD / %D	
	ICP interference check						
LCS	Laboratory control sample 1:53	Cr	38.783 ug/L	40.000 ug/L	97.0%	97.0%	Y
MS	Matrix spike 2:11	Cr	4.237 (SSR-SR) 41.031 - 4.237 = 36.794 ug/L	40.000 ug/L	92.0%	92.0%	Y
MSD	Duplicate	Cr	39.611 ug/L	Found: 41.021 ug/L	3.52%	3.52% RPD	Y
	ICP serial dilution						

Comments: Refer to appropriate worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 39385B4a

SDG #: 17-20934

VALIDATION FINDINGS WORKSHEET

Sample Calculation Verification

Page: 1 of 1

Reviewer: JB

2nd reviewer: KK

METHOD: Trace metals (EPA CLP SOW ILM02.1)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

Y N N/A Have results been reported and calculated correctly?

Y N N/A Are results within the calibrated range of the instruments and within the linear range of the ICP?

Y N N/A Are all detection limits below the CRDL?

Detected analyte results for Cr were recalculated and verified using the following equation:

Concentration = $\frac{(RD)(FV)(Dil)}{(In. Vol.)}$

Recalculation:

From Raw Data: 1.833 ug/L

- RD = Raw data concentration
- FV = Final volume (ml)
- In. Vol. = Initial volume (ml) or weight (G)
- Dil = Dilution factor

#	Sample ID	Analyte	Reported Concentration (<u>ug/L</u>)	Calculated Concentration (<u>ug/L</u>)	Acceptable (Y/N)
	<u>7</u>	<u>Cr</u>	<u>1.8</u>	<u>1.8</u>	<u>Y</u>

Note: _____

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017

LDC Report Date: September 20, 2017

Parameters: Wet Chemistry

Validation Level: Level III & IV

Laboratory: BC Laboratories, Inc.

Sample Delivery Group (SDG): 17-20934

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-4-3	1720934-02	Water	07/31/17
MW-4-2	1720934-03	Water	07/31/17
MW-4-1	1720934-04	Water	07/31/17
MW-21-5	1720934-05	Water	07/31/17
MW-21-4	1720934-06	Water	07/31/17
MW-21-3	1720934-07	Water	07/31/17
MW-21-2**	1720934-08**	Water	07/31/17
MW-21-1	1720934-09	Water	07/31/17
EB-6-073117	1720934-10	Water	07/31/17
SB-2-073117	1720934-11	Water	07/31/17
MW-21-3MS	1720934-07MS	Water	07/31/17
MW-21-3MSD	1720934-07MSD	Water	07/31/17
MW-21-3DUP	1720934-07DUP	Water	07/31/17

**Indicates sample underwent Level IV validation

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following methods:

Hexavalent Chromium by Environmental Protection Agency (EPA) SW 846 Method 7196

Perchlorate by EPA Method 314.0

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results. Samples appended with a double asterisk on the cover page were subjected to Level IV data validation, which is comprised of the QC summary forms as well as the raw data, to confirm sample quantitation and identification.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. Initial Calibration

All criteria for the initial calibration of each method were met.

III. Continuing Calibration

Continuing calibration frequency and analysis criteria were met for each method when applicable.

IV. Laboratory Blanks

Laboratory blanks were analyzed as required by the methods. No contaminants were found in the laboratory blanks.

V. Field Blanks

Sample EB-6-073117 was identified as an equipment blank. No contaminants were found with the following exceptions:

Blank ID	Analyte	Concentration
EB-6-073117	Perchlorate	0.75 ug/L

Sample SB-2-073117 was identified as a source blank. No contaminants were found.

VI. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

VIII. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the methods. Percent recoveries (%R) were within QC limits.

IX. Field Duplicates

No field duplicates were identified in this SDG.

X. Sample Result Verification

All sample result verifications were acceptable for samples which underwent Level IV validation. Raw data were not reviewed for Level III validation.

XI. Overall Assessment of Data

The analysis was conducted within all specifications of the methods. No results were rejected in this SDG.

The quality control criteria reviewed were met and are considered acceptable. Based upon the data validation all results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Wet Chemistry - Data Qualification Summary - SDG 17-20934

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 17-20934

No Sample Data Qualified in this SDG



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-4-3

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-02</u>	File ID: <u>F080917.seq-68.0000.txt</u>	
Sampled: <u>07/31/17 07:35</u>	Prepared: <u>08/09/17 12:00</u>	Analyzed: <u>08/10/17 13:08</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH0982</u>	Sequence: <u>1714391</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

172117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-4-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-03RE2</u>	File ID: <u>F080917.seq-98.0000.txt</u>	
Sampled: <u>07/31/17 08:05</u>	Prepared: <u>08/09/17 12:00</u>	Analyzed: <u>08/10/17 20:49</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH0982</u>	Sequence: <u>1714391</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	4.3	1		EPA-314.0

JDO92117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:59:28PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-314.0

MW-4-1

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-04

File ID: F080917.seq-71.0000.txt

Sampled: 07/31/17 08:30

Prepared: 08/09/17 12:00

Analyzed: 08/10/17 13:50

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H0982

Sequence: 1714391

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

209217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-21-5

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-05</u>	File ID: <u>F080917.seq-72.0000.txt</u>	
Sampled: <u>07/31/17 10:15</u>	Prepared: <u>08/09/17 12:00</u>	Analyzed: <u>08/10/17 14:03</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH0982</u>	Sequence: <u>1714391</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.3	1	J	EPA-314.0

5692117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-21-4

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-06</u>	File ID: <u>F080917.seq-74.0000.txt</u>	
Sampled: <u>07/31/17 09:45</u>	Prepared: <u>08/09/17 12:00</u>	Analyzed: <u>08/10/17 14:31</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH0982</u>	Sequence: <u>1714391</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.3	1	J	EPA-314.0

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:59:28PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-21-3

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-07

File ID: F080917.seq-76.0000.txt

Sampled: 07/31/17 10:50

Prepared: 08/09/17 12:00

Analyzed: 08/10/17 14:59

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0982

Sequence: 1714391

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	3.0	1	J	EPA-314.0

5609117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:59:28PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-314.0

MW-21-2

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-08

File ID: F080917.seq-83.0000.txt

Sampled: 07/31/17 11:30

Prepared: 08/09/17 12:00

Analyzed: 08/10/17 16:36

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H0982

Sequence: 1714391

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	1.1	1	J	EPA-314.0

8092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-21-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-09</u>
Sampled: <u>07/31/17 12:00</u>	Prepared: <u>08/09/17 12:00</u>
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>
Batch: <u>B[H0982</u>	Sequence: <u>1714391</u>
	Calibration: <u>UNASSIGNED</u>
	Instrument: <u>IC6</u>
	File ID: <u>F080917.seq-85.0000.txt</u>
	Analyzed: <u>08/10/17 17:03</u>
	Initial/Final: <u>20 ml / 20 ml</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	7.7	1		EPA-314.0

86092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

EB-6-073117

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-10</u>	File ID: <u>F080917.seq-87.0000.txt</u>	
Sampled: <u>07/31/17 12:15</u>	Prepared: <u>08/09/17 12:00</u>	Analyzed: <u>08/10/17 17:31</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BH0982</u>	Sequence: <u>1714391</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.75	1	J	EPA-314.0

82092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:59:28PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

SB-2-073117

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-11

File ID: F080917.seq-88.0000.txt

Sampled: 07/31/17 12:30

Prepared: 08/09/17 12:00

Analyzed: 08/10/17 17:45

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH0982

Sequence: 1714391

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

8/09/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:59:28PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-4-3

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-02

File ID: 170731 1957 CR6-035

Sampled: 07/31/17 07:35

Prepared: 07/31/17 19:57

Analyzed: 07/31/17 20:15

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0129

Sequence:

1713707

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

2/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-4-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-03</u>	File ID: <u>170731 1957 CR6-036</u>	
Sampled: <u>07/31/17 08:05</u>	Prepared: <u>07/31/17 19:57</u>	Analyzed: <u>07/31/17 20:15</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH0129</u>	Sequence: <u>1713707</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-4-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-04</u>	File ID: <u>170731 1957 CR6-037</u>	
Sampled: <u>07/31/17 08:30</u>	Prepared: <u>07/31/17 19:57</u>	Analyzed: <u>07/31/17 20:15</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH0129</u>	Sequence: <u>1713707</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/25/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-21-5

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-05

File ID: 170731 1957 CR6-038

Sampled: 07/31/17 10:15

Prepared: 07/31/17 19:57

Analyzed: 07/31/17 20:15

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0129

Sequence: 1713707

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0014	1	J	EPA-7196

209217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-21-4

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-06</u>	File ID: <u>170731 1957 CR6-065</u>	
Sampled: <u>07/31/17 09:45</u>	Prepared: <u>07/31/17 19:57</u>	Analyzed: <u>07/31/17 21:01</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH0129</u>	Sequence: <u>1713707</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0013	1	J	EPA-7196

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-21-3

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-07</u>	File ID: <u>170731 1957 CR6-027</u>	
Sampled: <u>07/31/17 10:50</u>	Prepared: <u>07/31/17 19:57</u>	Analyzed: <u>07/31/17 19:57</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH0129</u>	Sequence: <u>1713707</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/25/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-21-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-08</u>	File ID: <u>170731 1957 CR6-040</u>	
Sampled: <u>07/31/17 11:30</u>	Prepared: <u>07/31/17 19:57</u>	Analyzed: <u>07/31/17 20:15</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B[H0129</u>	Sequence: <u>1713707</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/25/2017 12:59:28PM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-21-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-20934</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1720934-09</u>	File ID: <u>170731 1957 CR6-074</u>	
Sampled: <u>07/31/17 12:00</u>	Prepared: <u>07/31/17 19:57</u>	Analyzed: <u>07/31/17 21:17</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BH0129</u>	Sequence: <u>1713707</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0018	1	J	EPA-7196

8/27/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:59:28PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

EB-6-073117

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-10

File ID: 170731 1957 CR6-042

Sampled: 07/31/17 12:15

Prepared: 07/31/17 19:57

Analyzed: 07/31/17 20:15

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0129

Sequence:

1713707

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

6/12/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/25/2017 12:59:28PM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

SB-2-073117

Laboratory: BC Laboratories

SDG: 17-20934

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1720934-11

File ID: 170731 1957 CR6-075

Sampled: 07/31/17 12:30

Prepared: 07/31/17 19:57

Analyzed: 07/31/17 21:17

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0129

Sequence:

1713707

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

82092117

LDC #: 39385B6

VALIDATION COMPLETENESS WORKSHEET

Date: 9/18/17

SDG #: 17-20934

Level III/IV

Page: 1 of 1

Laboratory: BC Laboratories, Inc.

Reviewer: V3

2nd Reviewer: V1

METHOD: (Analyte) Hexavalent Chromium (EPA SW846 Method 7196), Perchlorate (EPA Method 314.0)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A/A	
II	Initial calibration	A	
III.	Calibration verification	A	
IV	Laboratory Blanks	A	
V	Field blanks	SW	EB=9, SB=10 *
VI.	Matrix Spike/Matrix Spike Duplicates	A	
VII.	Duplicate sample analysis	A	
VIII.	Laboratory control samples	A	LCS
IX.	Field duplicates	N	
X.	Sample result verification	A	Not reviewed for Level III validation.
XI	Overall assessment of data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

* ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

** Indicates sample was underwent Level IV review

	Client ID	Lab ID	Matrix	Date
1	MW-4-3	1720934-02	Water	07/31/17
2	MW-4-2	1720934-03	Water	07/31/17
3	MW-4-1	1720934-04	Water	07/31/17
4	MW-21-5	1720934-05	Water	07/31/17
5	MW-21-4	1720934-06	Water	07/31/17
6	MW-21-3	1720934-07	Water	07/31/17
7	MW-21-2**	1720934-08**	Water	07/31/17
8	MW-21-1	1720934-09	Water	07/31/17
9	EB-6-073117	1720934-10	Water	07/31/17
10	SB-2-073117	1720934-11	Water	07/31/17
11	MW-21-3MS	1720934-07MS	Water	07/31/17
12	MW-21-3MSD	1720934-07MSD	Water	07/31/17
13	MW-21-3DUP	1720934-07DUP	Water	07/31/17
14				
15				
16				

Notes: _____

Method: Inorganics (EPA Method See Cover)

Validation Area	Yes	No	NA	Findings/Comments
<i>I. Technical holding times</i>				
All technical holding times were met.	✓			
<i>II. Calibration</i>				
Were all instruments calibrated daily, each set-up time?	✓			
Were the proper number of standards used?	✓			
Were all initial calibration correlation coefficients ≥ 0.995 ?	✓			
Were all initial and continuing calibration verification %Rs within the 90-110% QC limits?	✓			
Were titrant checks performed as required? (Level IV only)			✓	
Were balance checks performed as required? (Level IV only)			✓	
<i>III. Blanks</i>				
Was a method blank associated with every sample in this SDG?	✓			
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.		✓		
<i>IV. Matrix spike/Matrix spike duplicates and Duplicates</i>				
Were a matrix spike (MS) and duplicate (DUP) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD or MS/DUP. Soil / Water.	✓			
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the 75-125 QC limits? If the sample concentration exceeded the spike concentration by a factor of 4 or more, no action was taken.	✓			
Were the MS/MSD or duplicate relative percent differences (RPD) $\leq 20\%$ for waters and $\leq 35\%$ for soil samples? A control limit of \leq CRDL ($\leq 2X$ CRDL for soil) was used for samples that were $\leq 5X$ the CRDL, including when only one of the duplicate sample values were $\leq 5X$ the CRDL.	✓			
<i>V. Laboratory control samples</i>				
Was an LCS analyzed for this SDG?	✓			
Was an LCS analyzed per extraction batch?	✓			
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the 80-120% (85-115% for Method 300.0) QC limits?	✓			
<i>VI. Regional Quality Assurance and Quality Control</i>				
Were performance evaluation (PE) samples performed?			✓	
Were the performance evaluation (PE) samples within the acceptance limits?			✓	

VALIDATION FINDINGS CHECKLIST

Validation Area	Yes	No	NA	Findings/Comments
VII. Sample Result Verification				
Were RLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	/			
Were detection limits < RL?	/			
VIII. Overall assessment of data				
Overall assessment of data was found to be acceptable.	/			
IX. Field duplicates				
Field duplicate pairs were identified in this SDG.		/		
Target analytes were detected in the field duplicates.			/	
X. Field blanks				
Field blanks were identified in this SDG.	/			
Target analytes were detected in the field blanks.	/			

VALIDATION FINDINGS WORKSHEET
Sample Specific Analysis Reference

All circled methods are applicable to each sample.

Sample ID	Parameter
1-10	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
00	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
11-13	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄

Comments: _____

LDC #: 3938586
SDG #: 17-20934

VALIDATION FINDINGS WORKSHEET Field Blanks

Page: 1 of 1
Reviewer: [Signature]
2nd reviewer: [Signature]

METHOD: Inorganics

N N/A Were field blanks identified in this SDG?
 Y N N/A Were target analytes detected in the field blanks?

Sample: 9 Field Blank / Trip Blank / Rinsate / Other EB (circle one)

Analyte	Concentration Units ($\mu\text{g/L}$)
ClO_4	0.75

Sample: _____ Field Blank / Trip Blank / Rinsate / Other _____ (circle one)

Analyte	Concentration Units ()

LDC #: 39385B

**Validation Findings Worksheet
Initial and Continuing Calibration Calculation Verification**

Page: 1 of 1
 Reviewer: B
 2nd Reviewer: [Signature]

Method: Inorganics, Method See Cover

The correlation coefficient (r) for the calibration of ClO₄⁻ was recalculated. Calibration date: 8/11/17

An initial or continuing calibration verification percent recovery (%R) was recalculated for each type of analysis using the following formula:

$$\%R = \frac{\text{Found} \times 100}{\text{True}}$$

Where, Found = concentration of each analyte measured in the analysis of the ICV or CCV solution
 True = concentration of each analyte in the ICV or CCV source

Type of analysis	Analyte	Standard	Conc. (ug/L)	Area	Recalculated	Reported	Acceptable (Y/N)
					r or r ²	r or r ²	
Initial calibration	ClO ₄ ⁻	s1	0	0.0001	99.9421%	99.9141%	Y
		s2	2	0.0024			
		s3	4	0.0044			
		s4	6	0.007			
		s5	10	0.0119			
		s6	20	0.0243			
Calibration verification	Cr ⁶⁺	ICV	<u>Found:</u> 0.0498 mg/L	<u>True:</u> 0.050000 mg/L	100%	101%	Y
Calibration verification	ClO ₄ ⁻	CCV	<u>Found:</u> 9.135 ug/L	<u>True:</u> 10.000 ug/L	91.4%	93.9%	Y
Calibration verification							

Comments: Refer to Calibration Verification findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

VALIDATION FINDINGS WORKSHEET
Level IV Recalculation Worksheet

METHOD: Inorganics, Method See Cover

Percent recoveries (%R) for a laboratory control sample and a matrix spike sample were recalculated using the following formula:

$$\%R = \frac{\text{Found}}{\text{True}} \times 100$$
 Where, Found = concentration of each analyte measured in the analysis of the sample. For the matrix spike calculation, Found = SSR (spiked sample result) - SR (sample result).
 True = concentration of each analyte in the source.

A sample and duplicate relative percent difference (RPD) was recalculated using the following formula:

$$RPD = \frac{|S-D|}{(S+D)/2} \times 100$$
 Where, S = Original sample concentration
 D = Duplicate sample concentration

Sample ID	Type of Analysis	Element	Found / S (units)	True / D (units)	Recalculated	Reported	Acceptable (Y/N)
					%R / RPD	%R / RPD	
LCS	Laboratory control sample	ClO_4^-	9.9643 µg/L	10.000 µg/L	99.7%	97.2%	Y
MS	Matrix spike sample	Cr_6^+	ND (SSR-SR) 0.0498 = 0.05242 mg/L	0.052632 mg/L	99.6%	99.4%	Y
MSD	Duplicate sample	Cr_6^+	Found: 0.05242 mg/L	Found: 0.05242 mg/L	0% RPD	0.08667 RPD	Y

Comments: _____

VALIDATION FINDINGS WORKSHEET
Sample Calculation Verification

METHOD: Inorganics, Method See Cover

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

- Y N N/A Have results been reported and calculated correctly?
- Y N N/A Are results within the calibrated range of the instruments?
- Y N N/A Are all detection limits below the CRQL?

Compound (analyte) results for C104 reported with a positive detect were recalculated and verified using the following equation:

Concentration =

Recalculation:

$$\begin{aligned}
 y &= mx + b \\
 y &= 0.001 \\
 m &= 0.001203 \\
 b &= 4.7 E^{-6}
 \end{aligned}$$

$$C_{104} = \frac{0.001 - 4.7 E^{-6}}{0.001203} = 0.8269 \mu g/L$$

#	Sample ID	Analyte	Reported Concentration ($\mu g/L$)	Calculated Concentration ($\mu g/L$)	Acceptable (Y/N)
	<u>7</u>	<u>C104</u>	<u>1.1</u>	<u>0.83</u>	<u>Y</u>

Note: _____

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017
LDC Report Date: September 21, 2017
Parameters: Volatiles
Validation Level: Level III
Laboratory: BC Laboratories, Inc.
Sample Delivery Group (SDG): 17-21317

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
TB-8-080217	1721317-01	Water	08/02/17
MW-13	1721317-02	Water	08/02/17
DUP-6-3Q17	1721317-03	Water	08/02/17
MW-6	1721317-04	Water	08/02/17
DUP-7-3Q17	1721317-05	Water	08/02/17
MW-5	1721317-06	Water	08/02/17
MW-8	1721317-08	Water	08/02/17
EB-8-080217	1721317-09	Water	08/02/17
MW-8MS	1721317-08MS	Water	08/02/17
MW-8MSD	1721317-08MSD	Water	08/02/17

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Superfund Organic Methods Data Review (June 2008). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following method:

Volatile Organic Compounds (VOCs) by Environmental Protection Agency (EPA) Method 524.2

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition and cooler temperatures upon receipt met validation criteria.

All technical holding time requirements were met.

II. GC/MS Instrument Performance Check

Instrument performance was checked at 12 hour intervals.

All ion abundance requirements were met.

III. Initial Calibration and Initial Calibration Verification

An initial calibration was performed as required by the method.

For compounds where average relative response factors (RRFs) were utilized, the percent relative standard deviations (%RSD) were less than or equal to 20.0%.

In the case where the laboratory used a calibration curve to evaluate the compounds, all coefficients of determination (r^2) were greater than or equal to 0.990.

The percent differences (%D) of the initial calibration verification (ICV) standard were less than or equal to 30.0% for all compounds.

IV. Continuing Calibration

Continuing calibration was performed at the required frequencies.

The percent differences (%D) were less than or equal to 30.0% for all compounds with the following exceptions:

Date	Compound	%D	Associated Samples	Flag	A or P
08/07/17	Methyl iodide Pentachloroethane	52.3 (≤ 30) 32.8 (≤ 30)	TB-8-080217 MW-13 DUP-6-3Q17 MW-8	UJ (all non-detects) UJ (all non-detects)	P
08/08/17	Bromomethane Methyl iodide Pentachloroethane	32.6 (≤ 30) 52.1 (≤ 30) 36.2 (≤ 30)	MW-6 DUP-7-3Q17 MW-5 EB-8-080217	UJ (all non-detects) UJ (all non-detects) UJ (all non-detects)	P

V. Laboratory Blanks

Laboratory blanks were analyzed as required by the method. No contaminants were found in the laboratory blanks.

VI. Field Blanks

Sample TB-8-080217 was identified as a trip blank. No contaminants were found.

Sample EB-8-080217 was identified as an equipment blank. No contaminants were found.

VII. Surrogates

Surrogates were added to all samples as required by the method. All surrogate recoveries (%R) were within QC limits.

VIII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

IX. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the method. Percent recoveries (%R) were within QC limits.

X. Field Duplicates

Samples MW-13 and DUP-6-3Q17 and samples MW-6 and DUP-73Q17 were identified as field duplicates. No results were detected in any of the samples with the following exceptions:

Compound	Concentration (ug/L)		RPD (Limits)
	MW-13	DUP-6-3Q17	
Carbon tetrachloride	0.28	0.27	4
Chloroform	2.9	2.8	4
Tetrachloroethene	0.23	0.28	20
Trichloroethene	0.31	0.31	0

Compound	Concentration (ug/L)		RPD (Limits)
	MW-6	DUP-7-3Q17	
Chloroform	0.48	0.50	4
Tetrachloroethene	0.57	0.49	15
Trichloroethene	2.6	2.7	4

XI. Internal Standards

All internal standard areas and retention times were within QC limits.

XII. Compound Quantitation

Raw data were not reviewed for Level III validation.

XIII. Target Compound Identifications

Raw data were not reviewed for Level III validation.

XIV. System Performance

Raw data were not reviewed for Level III validation.

XV. Overall Assessment of Data

The analysis was conducted within all specifications of the method. No results were rejected in this SDG.

Due to continuing calibration %D, data were qualified as estimated in eight samples.

The quality control criteria reviewed, other than those discussed above, were met and are considered acceptable. Sample results that were found to be estimated (J) are usable for limited purposes only. Based upon the data validation all other results are considered valid and usable for all purposes.

NASA JPL, 2Q2017
Volatiles - Data Qualification Summary - SDG 17-21317

Sample	Compound	Flag	A or P	Reason
TB-8-080217 MW-13 DUP-6-3Q17 MW-8	Methyl iodide Pentachloroethane	UJ (all non-detects) UJ (all non-detects)	P	Continuing calibration (%D)
MW-6 DUP-7-3Q17 MW-5 EB-8-080217	Bromomethane Methyl iodide Pentachloroethane	UJ (all non-detects) UJ (all non-detects) UJ (all non-detects)	P	Continuing calibration (%D)

NASA JPL, 2Q2017
Volatiles - Laboratory Blank Data Qualification Summary - SDG 17-21317

No Sample Data Qualified in this SDG



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

TB-8-080217

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-01 File ID: 07AUG14.D
Sampled: 08/02/17 07:00 Prepared: 08/07/17 09:48 Analyzed: 08/07/17 16:57
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

Handwritten signature



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

TB-8-080217

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-01 File ID: 07AUG14.D
Sampled: 08/02/17 07:00 Prepared: 08/07/17 09:48 Analyzed: 08/07/17 16:57
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/30/17



Tidewater Inc. Reported: 8/30/2017 10:31:52AM
 3761 Attucks Drive Project: JPL- GW Monitoring Wells
 Powell, OH 43065 Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

TB-8-080217

Laboratory: BC Laboratories SDG: 17-21317
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1721317-01 File ID: 07AUG14.D
 Sampled: 08/02/17 07:00 Prepared: 08/07/17 09:48 Analyzed: 08/07/17 16:57
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UT</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UT</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	9.4700	94.7	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.7300	97.3	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.7600	97.6	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	231884	6.58	239780	6.57	
Chlorobenzene-d5 (IS)	92539	9.61	97357	9.62	
1,4-Difluorobenzene (IS)	349804	7.38	364372	7.38	

* Values outside of QC limits

809217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:31:52AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

TB-8-080217

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21317</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721317-01</u>	File ID:	<u>07AUG14.D</u>		
Sampled:	<u>08/02/17 07:00</u>	Prepared:	<u>08/07/17 09:48</u>	Analyzed:	<u>08/07/17 16:57</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0652</u>	Sequence:	<u>1713910</u>	Calibration:	<u>1708011</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

12/21/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:31:52AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-13

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21317</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721317-02</u>	File ID:	<u>07AUG15.D</u>		
Sampled:	<u>08/02/17 08:50</u>	Prepared:	<u>08/07/17 09:48</u>	Analyzed:	<u>08/07/17 17:20</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>B[H0652</u>	Sequence:	<u>1713910</u>	Calibration:	<u>1708011</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.28	J
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	2.9	
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

82 09-117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-13

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-02 File ID: 07AUG15.D
Sampled: 08/02/17 08:50 Prepared: 08/07/17 09:48 Analyzed: 08/07/17 17:20
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U UJ
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U UJ
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	9.9700	99.7	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.9700	99.7	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.6800	96.8	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	221801	6.57	239780	6.57	
Chlorobenzene-d5 (IS)	88784	9.62	97357	9.62	
1,4-Difluorobenzene (IS)	339157	7.38	364372	7.38	

* Values outside of QC limits

Handwritten number: 709217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:31:52AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-13

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21317</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721317-02</u>	File ID:	<u>07AUG15.D</u>		
Sampled:	<u>08/02/17 08:50</u>	Prepared:	<u>08/07/17 09:48</u>	Analyzed:	<u>08/07/17 17:20</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BJH0652</u>	Sequence:	<u>1713910</u>	Calibration:	<u>1708011</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

Handwritten: 10/9/2017



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

DUP-6-3Q17

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21317</u>
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721317-03</u>
		File ID:	<u>07AUG16.D</u>
Sampled:	<u>08/02/17 09:00</u>	Prepared:	<u>08/07/17 09:48</u>
		Analyzed:	<u>08/07/17 17:43</u>
Solids:		Preparation:	<u>EPA 5030 Water MS</u>
		Initial/Final:	<u>25 ml / 25 ml</u>
Batch:	<u>B[H0652</u>	Sequence:	<u>1713910</u>
		Calibration:	<u>1708011</u>
		Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.27	J
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	2.8	
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

DUP-6-3Q17

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-03 File ID: 07AUG16.D
Sampled: 08/02/17 09:00 Prepared: 08/07/17 09:48 Analyzed: 08/07/17 17:43
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.28	J
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.31	J
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

DUP-6-3Q17

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-03 File ID: 07AUG16.D
Sampled: 08/02/17 09:00 Prepared: 08/07/17 09:48 Analyzed: 08/07/17 17:43
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>VJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>VJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	9.5000	95.0	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.7800	97.8	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.6800	96.8	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	223060	6.57	239780	6.57	
Chlorobenzene-d5 (IS)	88076	9.62	97357	9.62	
1,4-Difluorobenzene (IS)	336925	7.38	364372	7.38	

* Values outside of QC limits

8/9/2017



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
 Project: JPL- GW Monitoring Wells
 Project Number: 3Q17
 Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
 EPA-524.2

MW-6

Laboratory: BC Laboratories SDG: 17-21317
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1721317-04 File ID: 07AUG40.D
 Sampled: 08/02/17 09:50 Prepared: 08/07/17 09:48 Analyzed: 08/08/17 02:55
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>UT</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.48	J
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-6

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-04 File ID: 07AUG40.D
Sampled: 08/02/17 09:50 Prepared: 08/07/17 09:48 Analyzed: 08/08/17 02:55
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,1,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.57	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	2.6	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/30/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-6

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-04 File ID: 07AUG40.D
Sampled: 08/02/17 09:50 Prepared: 08/07/17 09:48 Analyzed: 08/08/17 02:55
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.160	102	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.8400	98.4	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.3600	93.6	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	190606	6.57	194267	6.58	
Chlorobenzene-d5 (IS)	77660	9.61	74707	9.62	
1,4-Difluorobenzene (IS)	283916	7.38	288644	7.38	

* Values outside of QC limits

Handwritten: 7207247



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:31:52AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-6

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21317</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721317-04</u>	File ID:	<u>07AUG40.D</u>		
Sampled:	<u>08/02/17 09:50</u>	Prepared:	<u>08/07/17 09:48</u>	Analyzed:	<u>08/08/17 02:55</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>B/H0652</u>	Sequence:	<u>1713910</u>	Calibration:	<u>1708011</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

Handwritten signature/initials



Tidewater Inc. Reported: 8/30/2017 10:31:52AM
 3761 Attucks Drive Project: JPL- GW Monitoring Wells
 Powell, OH 43065 Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

DUP-7-3Q17

Laboratory: BC Laboratories SDG: 17-21317
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1721317-05 File ID: 07AUG41.D
 Sampled: 08/02/17 10:00 Prepared: 08/07/17 09:48 Analyzed: 08/08/17 03:18
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BJH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>U5</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.50	
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

DUP-7-3Q17

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-05 File ID: 07AUG41.D
Sampled: 08/02/17 10:00 Prepared: 08/07/17 09:48 Analyzed: 08/08/17 03:18
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.49	J
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	2.7	
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

1709217



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

DUP-7-3Q17

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-05 File ID: 07AUG41.D
Sampled: 08/02/17 10:00 Prepared: 08/07/17 09:48 Analyzed: 08/08/17 03:18
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>US</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>US</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.300	103	75 - 125	
Toluene-d8 (Surrogate)	10.000	10.230	102	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.6900	96.9	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	201103	6.57	194267	6.58	
Chlorobenzene-d5 (IS)	80695	9.61	74707	9.62	
1,4-Difluorobenzene (IS)	288704	7.38	288644	7.38	

* Values outside of QC limits

Handwritten signature

Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-5

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-06 File ID: 07AUG42.D
Sampled: 08/02/17 11:01 Prepared: 08/07/17 07:00 Analyzed: 08/08/17 03:41
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0523 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>WJ</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-5

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-06 File ID: 07AUG42.D
Sampled: 08/02/17 11:01 Prepared: 08/07/17 07:00 Analyzed: 08/08/17 03:41
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0523 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-5

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-06 File ID: 07AUG42.D
Sampled: 08/02/17 11:01 Prepared: 08/07/17 07:00 Analyzed: 08/08/17 03:41
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0523 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.500	105	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.6600	96.6	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.9600	99.6	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	187312	6.57	194267	6.58	
Chlorobenzene-d5 (IS)	71675	9.62	74707	9.62	
1,4-Difluorobenzene (IS)	276661	7.38	288644	7.38	

* Values outside of QC limits

8/30/17



<p>Tidewater Inc. 3761 Attucks Drive Powell, OH 43065</p>	<p>Reported: 8/30/2017 10:31:52AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner</p>
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-5

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-06</u>	File ID: <u>07AUG42.D</u>	
Sampled: <u>08/02/17 11:01</u>	Prepared: <u>08/07/17 07:00</u>	Analyzed: <u>08/08/17 03:41</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>B[H0523</u>	Sequence: <u>1713910</u>	Calibration: <u>1708011</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8092417



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-8

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-08 File ID: 07AUG07.D
Sampled: 08/02/17 13:01 Prepared: 08/07/17 07:00 Analyzed: 08/07/17 14:16
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0523 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/30/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-8

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-08 File ID: 07AUG07.D
Sampled: 08/02/17 13:01 Prepared: 08/07/17 07:00 Analyzed: 08/07/17 14:16
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0523 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,1,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	J
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

SLC 8/21/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-8

Laboratory: BC Laboratories SDG: 17-21317
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1721317-08 File ID: 07AUG07.D
 Sampled: 08/02/17 13:01 Prepared: 08/07/17 07:00 Analyzed: 08/07/17 14:16
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0523 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U US
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U US
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	8.9400	89.4	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.6800	96.8	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.7400	97.4	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	232617	6.57	239780	6.57	
Chlorobenzene-d5 (IS)	91169	9.62	97357	9.62	
1,4-Difluorobenzene (IS)	348181	7.38	364372	7.38	

* Values outside of QC limits

8692117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:31:52AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-8

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-08</u>	File ID: <u>07AUG07.D</u>	
Sampled: <u>08/02/17 13:01</u>	Prepared: <u>08/07/17 07:00</u>	Analyzed: <u>08/07/17 14:16</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BIH0523</u>	Sequence: <u>1713910</u>	Calibration: <u>1708011</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/09/2017



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-8-080217

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-09 File ID: 07AUG43.D
Sampled: 08/02/17 13:15 Prepared: 08/07/17 07:00 Analyzed: 08/08/17 04:04
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0523 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>W5</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/30/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-8-080217

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-09 File ID: 07AUG43.D
Sampled: 08/02/17 13:15 Prepared: 08/07/17 07:00 Analyzed: 08/08/17 04:04
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0523 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

Scanned



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-8-080217

Laboratory: BC Laboratories SDG: 17-21317
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721317-09 File ID: 07AUG43.D
Sampled: 08/02/17 13:15 Prepared: 08/07/17 07:00 Analyzed: 08/08/17 04:04
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0523 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.180	102	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.8400	98.4	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.8600	98.6	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	191839	6.58	194267	6.58	
Chlorobenzene-d5 (IS)	74543	9.61	74707	9.62	
1,4-Difluorobenzene (IS)	287431	7.38	288644	7.38	

* Values outside of QC limits

8/09/17



BC *Laboratories, Inc.*

Environmental Testing Laboratory Since 1949



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:31:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

EB-8-080217

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21317</u>
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721317-09</u>
Sampled:	<u>08/02/17 13:15</u>	Prepared:	<u>08/07/17 07:00</u>
Solids:		Preparation:	<u>EPA 5030 Water MS</u>
Batch:	<u>BH0523</u>	Sequence:	<u>1713910</u>
		Calibration:	<u>1708011</u>
		Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

LDC #: 39385C1
 SDG #: 17-21317
 Laboratory: BC Laboratories, Inc.

VALIDATION COMPLETENESS WORKSHEET
 Level III

Date: 9/2/17
 Page: 1 of 1
 Reviewer: [Signature]
 2nd Reviewer: YU

METHOD: GC/MS Volatiles (EPA Method 524.2)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A	
II.	GC/MS Instrument performance check	A	
III.	Initial calibration/ICV	A/A	RSD ≤ 20%. Y ² ICV = 33%
IV.	Continuing calibration	M	CCV ≤ 30%
V.	Laboratory Blanks	A	
VI.	Field blanks	ND	TB=1. EB=8
VII.	Surrogate spikes	A	
VIII.	Matrix spike/Matrix spike duplicates	A	
IX.	Laboratory control samples	A	ICS
X.	Field duplicates	M	D=2+3. A+5
XI.	Internal standards	A	
XII.	Compound quantitation RL/LOQ/LODs	N	
XIII.	Target compound identification	N	
XIV.	System performance	N	
XV.	Overall assessment of data	A	

Note: A = Acceptable
 N = Not provided/applicable
 SW = See worksheet
 ND = No compounds detected
 R = Rinsate
 FB = Field blank
 D = Duplicate
 TB = Trip blank
 EB = Equipment blank
 SB=Source blank
 OTHER:

	Client ID	Lab ID	Matrix	Date
1	TB-8-080217	1721317-01	Water	08/02/17
2	MW-13	1721317-02	Water	08/02/17
3	DUP-6-3Q17	1721317-03	Water	08/02/17
4	MW-6	1721317-04	Water	08/02/17
5	DUP-7-3Q17	1721317-05	Water	08/02/17
6	MW-5	1721317-06	Water	08/02/17
7	MW-8	1721317-08	Water	08/02/17
8	EB-8-080217	1721317-09	Water	08/02/17
9	MW-8MS	1721317-08MS	Water	08/02/17
10	MW-8MSD	1721317-08MSD	Water	08/02/17
11				
12				
13				

TARGET COMPOUND WORKSHEET

METHOD: VOA

A. Chloromethane	AA. Tetrachloroethene	AAA. 1,3,5-Trimethylbenzene	AAAA. Ethyl tert-butyl ether	A1. 1,3-Butadiene
B. Bromomethane	BB. 1,1,2,2-Tetrachloroethane	BBB. 4-Chlorotoluene	BBBB. tert-Amyl methyl ether	B1. Hexane
C. Vinyl chloride	CC. Toluene	CCC. tert-Butylbenzene	CCCC. 1-Chlorohexane	C1. Heptane
D. Chloroethane	DD. Chlorobenzene	DDD. 1,2,4-Trimethylbenzene	DDDD. Isopropyl alcohol	D1. Propylene
E. Methylene chloride	EE. Ethylbenzene	EEE. sec-Butylbenzene	EEEE. Acetonitrile	E1. Freon 11
F. Acetone	FF. Styrene	FFF. 1,3-Dichlorobenzene	FFFF. Acrolein	F1. Freon 12
G. Carbon disulfide	GG. Xylenes, total	GGG. p-Isopropyltoluene	GGGG. Acrylonitrile	G1. Freon 113
H. 1,1-Dichloroethene	HH. Vinyl acetate	HHH. 1,4-Dichlorobenzene	HHHH. 1,4-Dioxane	H1. Freon 114
I. 1,1-Dichloroethane	II. 2-Chloroethylvinyl ether	III. n-Butylbenzene	IIII. Isobutyl alcohol	I1. 2-Nitropropane
J. 1,2-Dichloroethene, total	JJ. Dichlorodifluoromethane	JJJ. 1,2-Dichlorobenzene	JJJJ. Methacrylonitrile	J1. Dimethyl disulfide
K. Chloroform	KK. Trichlorofluoromethane	KKK. 1,2,4-Trichlorobenzene	KKKK. Propionitrile	K1. 2,3-Dimethyl pentane
L. 1,2-Dichloroethane	LL. Methyl-tert-butyl ether	LLL. Hexachlorobutadiene	LLLL. Ethyl ether	L1. 2,4-Dimethyl pentane
M. 2-Butanone	MM. 1,2-Dibromo-3-chloropropane	MMM. Naphthalene	MMMM. Benzyl chloride	M1. 3,3-Dimethyl pentane
N. 1,1,1-Trichloroethane	NN. Methyl ethyl ketone	NNN. 1,2,3-Trichlorobenzene	NNNN. Iodomethane	N1. 2-Methylpentane
O. Carbon tetrachloride	OO. 2,2-Dichloropropane	OOO. 1,3,5-Trichlorobenzene	OOOO. 1,1-Difluoroethane	O1. 3-Methylpentane
P. Bromodichloromethane	PP. Bromochloromethane	PPP. trans-1,2-Dichloroethene	PPPP. Tetrahydrofuran	P1. 3-Ethylpentane
Q. 1,2-Dichloropropane	QQ. 1,1-Dichloropropene	QQQ. cis-1,2-Dichloroethene	QQQQ. Methyl acetate	Q1. 2,2-Dimethylpentane
R. cis-1,3-Dichloropropene	RR. Dibromomethane	RRR. m,p-Xylenes	RRRR. Ethyl acetate	R1. 2,2,3- Trimethylbutane
S. Trichloroethene	SS. 1,3-Dichloropropane	SSS. o-Xylene	SSSS. Cyclohexane	S1. 2,2,4-Trimethylpentane
T. Dibromochloromethane	TT. 1,2-Dibromoethane	TTT. 1,1,2-Trichloro-1,2,2-trifluoroethane	TTTT. Methylcyclohexane	T1. 2-Methylhexane
U. 1,1,2-Trichloroethane	UU. 1,1,1,2-Tetrachloroethane	UUU. 1,2-Dichlorotetrafluoroethane	UUUU. Allyl chloride	U1. Nonanal
V. Benzene	VV. Isopropylbenzene	VVV. 4-Ethyltoluene	VVVV. Methyl methacrylate	V1. 2-Methylnaphthalene
W. trans-1,3-Dichloropropene	WW. Bromobenzene	WWW. Ethanol	WWWWW. Ethyl methacrylate	W1. Methanol
X. Bromoform	XX. 1,2,3-Trichloropropane	XXX. Di-isopropyl ether	XXXX. cis-1,4-Dichloro-2-butene	X1. 1,2,3-Trimethylbenzene
Y. 4-Methyl-2-pentanone	YY. n-Propylbenzene	YYY. tert-Butanol	YYYY. trans-1,4-Dichloro-2-butene	Y1.
Z. 2-Hexanone	ZZ. 2-Chlorotoluene	ZZZ. tert-Butyl alcohol	ZZZZ. Pentachloroethane	Z1.

LDC #: 39385c1

VALIDATION FINDINGS WORKSHEET
Continuing Calibration

Reviewer: Q
2nd Reviewer: KR

METHOD: GC/MS VOA (EPA Method 524.2)

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

Y N N/A Was a continuing calibration standard analyzed at least once every 12 hours for each instrument?

Y N N/A Were all percent differences (%D) < 30%?

#	Date	Standard ID	Compound	Finding %D (Limit: <30.0%)	Associated Samples	Qualifications
	8/7/17	07AUG03	Methyl iodide 2222	52.3 32.8	1-3.7.9-10. MB (ND)	✓ N/A ↓
	8/8/17	07AUG33	B Methyl iodide 2222	32.6 52.1 36.2	4-6.8.MB (ND)	✓ N/A ↓

LDC#: 39385C1

VALIDATION FINDINGS WORKSHEET
Field Duplicates

Page: 1 of 1
Reviewer: [Signature]
2nd Reviewer: [Signature]

METHOD: GCMS VOA (EPA Method 524.2)

Compound	Concentration (ug/L)		RPD
	2	3	
O	0.28	0.27	4
K	2.9	2.8	4
AA	0.23	0.28	20
S	0.31	0.31	0

Compound	Concentration (ug/L)		RPD
	4	5	
K	0.48	0.50	4
AA	0.57	0.49	15
S	2.6	2.7	4

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017
LDC Report Date: September 21, 2017
Parameters: Chromium
Validation Level: Level III
Laboratory: BC Laboratories, Inc.
Sample Delivery Group (SDG): 17-21317

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-13	1721317-02	Water	08/02/17
DUP-6-3Q17	1721317-03	Water	08/02/17
MW-6	1721317-04	Water	08/02/17
DUP-7-3Q17	1721317-05	Water	08/02/17
MW-5	1721317-06	Water	08/02/17
MW-15	1721317-07	Water	08/02/17
MW-8	1721317-08	Water	08/02/17
EB-8-080217	1721317-09	Water	08/02/17
MW-15MS	1721317-07MS	Water	08/02/17
MW-15MSD	1721317-07MSD	Water	08/02/17
MW-15DUP	1721317-07DUP	Water	08/02/17

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following method:

Chromium by Environmental Protection Agency (EPA) Method 200.8

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. ICPMS Tune

The mass calibration was within 0.1 AMU and the percent relative standard deviation (%RSD) was less than or equal to 5%.

III. Instrument Calibration

Initial and continuing calibrations were performed as required by the methods.

The initial calibration verification (ICV) and continuing calibration verification (CCV) standards were within QC limits.

IV. ICP Interference Check Sample Analysis

ICP interference check sample analyses were not required by the method.

V. Laboratory Blanks

Laboratory blanks were analyzed as required by the method. No contaminants were found in the laboratory blanks with the following exceptions:

Blank ID	Analyte	Maximum Concentration	Associated Samples
PB (prep blank)	Chromium	0.52500 ug/L	All samples in SDG 17-21317

Data qualification by the laboratory blanks was based on the maximum contaminant concentration in the laboratory blanks in the analysis of each analyte. The sample concentrations were either not detected or were significantly greater (>5X blank contaminants) than the concentrations found in the associated laboratory blanks with the following exceptions:

Sample	Analyte	Reported Concentration	Modified Final Concentration
MW-5	Chromium	1.2 ug/L	1.2U ug/L
MW-8	Chromium	1.9 ug/L	1.9U ug/L
EB-8-080217	Chromium	1.1 ug/L	1.1U ug/L

VI. Field Blanks

Sample EB-8-080217 was identified as an equipment blank. No contaminants were found with the following exceptions:

Blank ID	Analyte	Concentration (ug/L)
EB-8-080217	Chromium	1.1 ug/L

VII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VIII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

IX. Serial Dilution

Serial dilution was not performed for this SDG.

X. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the method. Percent recoveries (%R) were within QC limits.

XI. Field Duplicates

Samples MW-13 and DUP-6-3Q17 and samples MW-6 and DUP-7-3Q17 were identified as field duplicates. No results were detected in any of the samples with the following exceptions:

Analyte	Concentration (ug/L)		RPD
	MW-13	DUP-6-3Q17	
Chromium	1300	680	63

Analyte	Concentration (ug/L)		RPD
	MW-6	DUP-7-3Q17	
Chromium	30	120	120

XII. Internal Standards (ICP-MS)

Raw data were not reviewed for Level III validation.

XIII. Sample Result Verification

Raw data were not reviewed for Level III validation.

XIV. Overall Assessment of Data

The analysis was conducted within all specifications of the method. No results were rejected in this SDG.

Due to laboratory blank contamination, data were qualified as not detected in three samples.

The quality control criteria reviewed, other than those discussed above, were met and are considered acceptable. Based upon the data validation all other results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Metals - Data Qualification Summary - SDG 17-21317

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017
Metals - Laboratory Blank Data Qualification Summary - SDG 17-21317

Sample	Analyte	Modified Final Concentration	A or P
MW-5	Chromium	1.2U ug/L	A
MW-8	Chromium	1.9U ug/L	A
EB-8-080217	Chromium	1.1U ug/L	A



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:37:01AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-13

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-02

File ID: PE-EL3 170808-214

Sampled: 08/02/17 08:50

Prepared: 08/04/17 09:00

Analyzed: 08/09/17 02:19

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0425

Sequence: 1714081

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1300	1		EPA-200.8

U9217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:37:01AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-200.8

DUP-6-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-03</u>
Sampled: <u>08/02/17 09:00</u>	Prepared: <u>08/04/17 09:00</u>
Solids: <u>0.00</u>	Preparation: <u>EPA 200.2</u>
Batch: <u>BIH0425</u>	Sequence: <u>1714081</u>
	Calibration: <u>UNASSIGNED</u>
	Instrument: <u>PE-EL3</u>
	File ID: <u>PE-EL3_170808-215</u>
	Analyzed: <u>08/09/17 02:23</u>
	Initial/Final: <u>50 ml / 50 ml</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	680	1		EPA-200.8

1089217



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:37:01AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-6

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-04

File ID: PE-EL3_170808-216

Sampled: 08/02/17 09:50

Prepared: 08/04/17 09:00

Analyzed: 08/09/17 02:26

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0425

Sequence: 1714081

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	30	1		EPA-200.8

7609-117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:37:01AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

DUP-7-3Q17

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-05

File ID: PE-EL3 170808-217

Sampled: 08/02/17 10:00

Prepared: 08/04/17 09:00

Analyzed: 08/09/17 02:30

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0425

Sequence: 1714081

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	120	1		EPA-200.8

1209-117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:37:01AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-5

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-06</u>	File ID: <u>PE-EL3 170808-218</u>	
Sampled: <u>08/02/17 11:01</u>	Prepared: <u>08/04/17 09:00</u>	Analyzed: <u>08/09/17 02:33</u>	
Solids: <u>0.00</u>	Preparation: <u>EPA 200.2</u>	Initial/Final: <u>50 ml / 50 ml</u>	
Batch: <u>BJH0425</u>	Sequence: <u>1714081</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>PE-EL3</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.2 <i>u</i>	1	J	EPA-200.8

10/2/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:37:01AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-15

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-07</u>	File ID: <u>PE-EL3 170808-204</u>	
Sampled: <u>08/02/17 11:50</u>	Prepared: <u>08/04/17 09:00</u>	Analyzed: <u>08/09/17 01:45</u>	
Solids: <u>0.00</u>	Preparation: <u>EPA 200.2</u>	Initial/Final: <u>50 ml / 50 ml</u>	
Batch: <u>BIH0425</u>	Sequence: <u>1714081</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>PE-EL3</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	5.9	1		EPA-200.8

SLG 8/21/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:37:01AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-8

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-08

File ID: PE-EL3_170808-219

Sampled: 08/02/17 13:01

Prepared: 08/04/17 09:00

Analyzed: 08/09/17 02:37

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0425

Sequence: 1714081

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.9 U	1	J	EPA-200.8

8/09/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:37:01AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

EB-8-080217

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-09

File ID: PE-EL3 170808-220

Sampled: 08/02/17 13:15

Prepared: 08/04/17 09:00

Analyzed: 08/09/17 02:40

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0425

Sequence: 1714081

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.1 <i>u</i>	1	J	EPA-200.8

8/30/17

LDC #: 39385C4a

VALIDATION COMPLETENESS WORKSHEET

Date: 9/13/17

SDG #: 17-21317

Level III

Page: 1 of 1

Laboratory: BC Laboratories, Inc.

Reviewer: JB

2nd Reviewer: JT

METHOD: Chromium (EPA Method 200.8)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A/A	
II.	ICP/MS Tune	A	
III.	Instrument Calibration	A	
IV.	ICP Interference Check Sample (ICS) Analysis	N	Not Required
V.	Laboratory Blanks	SW	
VI.	Field Blanks	SW	EB=8
VII.	Matrix Spike/Matrix Spike Duplicates	A	
VIII.	Duplicate sample analysis	A	
IX.	Serial Dilution	N	Not performed
X.	Laboratory control samples	A	LCS
XI.	Field Duplicates	SW	(1,2) (3,4)
XII.	Internal Standard (ICP-MS)	N	Not Reviewed for level III
XIII.	Sample Result Verification	N	
XIV.	Overall Assessment of Data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	MW-13	1721317-02	Water	08/02/17
2	DUP-6-3Q17	1721317-03	Water	08/02/17
3	MW-6	1721317-04	Water	08/02/17
4	DUP-7-3Q17	1721317-05	Water	08/02/17
5	MW-5	1721317-06	Water	08/02/17
6	MW-15	1721317-07	Water	08/02/17
7	MW-8	1721317-08	Water	08/02/17
8	EB-8-080217	1721317-09	Water	08/02/17
9	MW-15MS	1721317-07MS	Water	08/02/17
10	MW-15MSD	1721317-07MSD	Water	08/02/17
11	MW-15DUP	1721317-07DUP	Water	08/02/17
12				
13				
14				

Notes:

VALIDATION FINDINGS WORKSHEET
PB/ICB/CCB QUALIFIED SAMPLES

METHOD: Trace metals (EPA SW 864 Method 6010B/6020/7000)
 Sample Concentration units, unless otherwise noted: ug/L

Soil preparation factor applied: NA
 Associated Samples: All

Analyte	Maximum PB ^a (mg/Kg)	Maximum PB ^a (ug/L)	Maximum ICB/CCB ^a (mg/L)	Action Level	5	7	8						
Cr		0.52500		2.625	1.2	1.9	1.1						

Samples with analyte concentrations within five times the associated ICB, CCB or PB concentration are listed above with the identifications from the Validation Completeness Worksheet. These sample results were qualified as not detected, "U".

Note : a - The listed analyte concentration is the highest ICB, CCB, or PB detected in the analysis of each element.

LDC #: 39385C4a
SDG #: 17-21317

VALIDATION FINDINGS WORKSHEET

Field Blanks

Page: 1 of 1
Reviewer: VB
2nd reviewer: VK

METHOD: Trace Metals (EPA CLP SOW ILM02.1)

Y N N/A Were field blanks identified in this SDG?
 Y N N/A Were target analytes detected in the field blanks?

Sample: 8 Field Blank / Trip Blank / Rinsate / Other EB (circle one)

Analyte	Concentration Units (<u>ug/L</u>)
<u>Cr</u>	<u>1.1</u>

Sample: _____ Field Blank / Trip Blank / Rinsate / Other _____ (circle one)

Analyte	Concentration Units ()

LDC#: 39385C4a

VALIDATION FINDINGS WORKSHEET
Field Duplicates

Page: 1 of 1
Reviewer: J3
2nd Reviewer: ST

METHOD: Metals (EPA Method 200.8)

Analyte	Concentration (ug/L)		RPD	
	1	2		
Chromium	1300	680	63	

Analyte	Concentration (ug/L)		RPD	
	3	4		
Chromium	30	120	120	

V:\FIELD DUPLICATES\Field Duplicates\FD_inorganic\2017\39385C4a.wpd

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017
LDC Report Date: September 21, 2017
Parameters: Wet Chemistry
Validation Level: Level III & IV
Laboratory: BC Laboratories, Inc.
Sample Delivery Group (SDG): 17-21317

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-13	1721317-02	Water	08/02/17
DUP-6-3Q17	1721317-03	Water	08/02/17
MW-6	1721317-04	Water	08/02/17
DUP-7-3Q17	1721317-05	Water	08/02/17
MW-5	1721317-06	Water	08/02/17
MW-15	1721317-07	Water	08/02/17
MW-8	1721317-08	Water	08/02/17
EB-8-080217	1721317-09	Water	08/02/17
MW-5MS	1721317-06MS	Water	08/02/17
MW-5MSD	1721317-06MSD	Water	08/02/17
MW-5DUP	1721317-06DUP	Water	08/02/17
MW-15MS	1721317-07MS	Water	08/02/17
MW-15MSD	1721317-07MSD	Water	08/02/17
MW-15DUP	1721317-07DUP	Water	08/02/17
MW-8MS	1721317-08MS	Water	08/02/17
MW-8MSD	1721317-08MSD	Water	08/02/17
MW-8DUP	1721317-08DUP	Water	08/02/17

**Indicates sample underwent Level IV validation

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following methods:

Chloride, Nitrate as Nitrogen, and Sulfate by Environmental Protection Agency (EPA) Method 300.0

Nitrite as Nitrogen by EPA 353.2

Hexavalent Chromium by EPA SW 846 Method 7196

Orthophosphate as Phosphorus by EPA Method 365.1

Perchlorate by EPA Method 314.0

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results. Samples appended with a double asterisk on the cover page were subjected to Level IV data validation, which is comprised of the QC summary forms as well as the raw data, to confirm sample quantitation and identification.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. Initial Calibration

All criteria for the initial calibration of each method were met.

III. Continuing Calibration

Continuing calibration frequency and analysis criteria were met for each method when applicable.

IV. Laboratory Blanks

Laboratory blanks were analyzed as required by the methods. No contaminants were found in the laboratory blanks with the following exceptions:

Blank ID	Analyte	Maximum Concentration	Associated Samples
PB (prep blank)	Nitrite as Nitrogen	0.012280 mg/L	MW-13 MW-8
ICB/CCB	Nitrite as Nitrogen	0.012586 mg/L	MW-13 MW-8

Data qualification by the laboratory blanks was based on the maximum contaminant concentration in the laboratory blanks in the analysis of each analyte. The sample concentrations were either not detected or were significantly greater (>5X blank contaminants) than the concentrations found in the associated laboratory blanks with the following exceptions:

Sample	Analyte	Reported Concentration	Modified Final Concentration
MW-13	Nitrite as Nitrogen	0.015 mg/L	0.015U mg/L
MW-8	Nitrite as Nitrogen	0.014 mg/L	0.014U mg/L

V. Field Blanks

Sample EB-8-080217 was identified as an equipment blank. No contaminants were found.

VI. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

VIII. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the methods. Percent recoveries (%R) were within QC limits.

IX. Field Duplicates

Samples MW-13 and DUP-6-3Q17 and samples MW-6 and DUP-7-3Q17 were identified as field duplicates. No results were detected in any of the samples with the following exceptions:

Analyte	Concentration		RPD
	MW-13	DUP-6-2Q17	
Perchlorate	220 ug/L	230 ug/L	4
Hexavalent chromium	0.0025 mg/L	0.0027 mg/L	8

Analyte	Concentration		RPD
	MW-6	DUP-7-2Q17	
Perchlorate	2.7 ug/L	2.3 ug/L	16
Hexavalent chromium	0.00070U mg/L	0.0020 mg/L	96

X. Sample Result Verification

All sample result verifications were acceptable for samples which underwent Level IV validation. Raw data were not reviewed for Level III validation.

XI. Overall Assessment of Data

The analysis was conducted within all specifications of the methods. No results were rejected in this SDG.

Due to laboratory blank contamination, data were qualified as not detected in two samples.

The quality control criteria reviewed, other than those discussed above, were met and are considered acceptable. Based upon the data validation all other results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Wet Chemistry - Data Qualification Summary - SDG 17-21317

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 17-21317

Sample	Analyte	Modified Final Concentration	A or P
MW-13	Nitrite as Nitrogen	0.015U mg/L	A
MW-8	Nitrite as Nitrogen	0.014U mg/L	A



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-300.0

MW-13

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-02

File ID: E080217A.seq-22

Sampled: 08/02/17 08:50

Prepared: 08/02/17 22:00

Analyzed: 08/03/17 04:39

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0315

Sequence: 1714191

Calibration: UNASSIGNED

Instrument: IC5

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
16887-00-6	Chloride	53	1		EPA-300.0
14797-55-8	Nitrate as N	4.8	1		EPA-300.0
14808-79-8	Sulfate	49	1		EPA-300.0

1092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:34:03AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-300.0

MW-8

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-08</u>	File ID: <u>E080217A.seq-23</u>	
Sampled: <u>08/02/17 13:01</u>	Prepared: <u>08/02/17 22:00</u>	Analyzed: <u>08/03/17 04:57</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B[H0315</u>	Sequence: <u>1714191</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC5</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
16887-00-6	Chloride	14	1		EPA-300.0
14797-55-8	Nitrate as N	2.6	1		EPA-300.0
14808-79-8	Sulfate	42	1		EPA-300.0

12/09/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-13

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-02RE2

File ID: F081317.seq-38.0000.txt

Sampled: 08/02/17 08:50

Prepared: 08/13/17 11:00

Analyzed: 08/14/17 08:25

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH1278

Sequence: 1714620

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	220	20	D	EPA-314.0

809217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:34:03AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

DUP-6-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-03RE2</u>	File ID: <u>F081317.seq-39.0000.txt</u>	
Sampled: <u>08/02/17 09:00</u>	Prepared: <u>08/13/17 11:00</u>	Analyzed: <u>08/14/17 08:39</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH1278</u>	Sequence: <u>1714620</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	230	20	D	EPA-314.0

82092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:34:03AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-314.0

MW-6

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-04RE4

File ID: F081617.seq-42.0000.txt

Sampled: 08/02/17 09:50

Prepared: 08/13/17 11:00

Analyzed: 08/16/17 11:37

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H1593

Sequence: 1714795

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.7	1	J	EPA-314.0

8/30/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:34:03AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
--	--

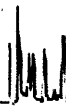
INORGANIC ANALYSIS DATA SHEET
EPA-314.0

DUP-7-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-05RE4</u>	File ID: <u>F081617.seq-43.0000.txt</u>	
Sampled: <u>08/02/17 10:00</u>	Prepared: <u>08/13/17 11:00</u>	Analyzed: <u>08/16/17 11:51</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BH1593</u>	Sequence: <u>1714795</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.3	1	J	EPA-314.0

8/30/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:34:03AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-5

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-06</u>	File ID: <u>F081317.seq-22.0000.txt</u>	
Sampled: <u>08/02/17 11:01</u>	Prepared: <u>08/13/17 11:00</u>	Analyzed: <u>08/13/17 15:56</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH1278</u>	Sequence: <u>1714620</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

807217



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-8

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-08RE3

File ID: F081317.seq-47.0000.txt

Sampled: 08/02/17 13:01

Prepared: 08/13/17 11:00

Analyzed: 08/14/17 11:01

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH1278

Sequence:

1714620

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

8/30/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

EB-8-080217

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-09RE2

File ID: F081317.seq-42.0000.txt

Sampled: 08/02/17 13:15

Prepared: 08/13/17 11:00

Analyzed: 08/14/17 09:20

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH1278

Sequence:

1714620

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

Handwritten signature/initials



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-353.2

MW-13

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-02

File ID: 170803 0842 NO2-032

Sampled: 08/02/17 08:50

Prepared: 08/03/17 08:42

Analyzed: 08/03/17 08:42

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0398

Sequence: 1713831

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
14797-65-0	Nitrite as N	0.015 U	1	J	EPA-353.2

8/30/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:34:03AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-353.2

MW-8

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-08

File ID: 170803 0842 NO2-028

Sampled: 08/02/17 13:01

Prepared: 08/03/17 08:42

Analyzed: 08/03/17 08:42

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B/H0398

Sequence: 1713831

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
14797-65-0	Nitrite as N	0.014 U	1	J	EPA-353.2

8657217



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-365.1

MW-13

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-02

File ID: 170803 0956 PO4-095

Sampled: 08/02/17 08:50

Prepared: 08/03/17 09:56

Analyzed: 08/03/17 09:57

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0397

Sequence: 1713938

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
---	ortho-Phosphate as P	0.044	1	J	EPA-365.1

8/30/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-365.1

MW-8

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-08

File ID: 170803 0956 PO4-091

Sampled: 08/02/17 13:01

Prepared: 08/03/17 09:56

Analyzed: 08/03/17 09:57

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0397

Sequence: 1713938

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
---	ortho-Phosphate as P	0.017	1	U	EPA-365.1

8/30/17

Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-13

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-02

File ID: 170802 2014 CR6-027

Sampled: 08/02/17 08:50

Prepared: 08/02/17 20:14

Analyzed: 08/02/17 20:32

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BJH0293

Sequence:

1713798

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0025	1		EPA-7196

8/30/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:34:03AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

DUP-6-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-03</u>
Sampled: <u>08/02/17 09:00</u>	File ID: <u>170802 2014 CR6-006</u>
Solids: <u>0.00</u>	Prepared: <u>08/02/17 20:14</u>
Batch: <u>B[H0293</u>	Analyzed: <u>08/02/17 20:14</u>
Sequence: <u>1713798</u>	Initial/Final: <u>20 ml / 20 ml</u>
	Preparation: <u>No Prep</u>
	Calibration: <u>UNASSIGNED</u>
	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0027	1		EPA-7196

8/30/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-7196

MW-6

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-04

File ID: 170802 2014 CR6-007

Sampled: 08/02/17 09:50

Prepared: 08/02/17 20:14

Analyzed: 08/02/17 20:14

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0293

Sequence: 1713798

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/30/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-7196

DUP-7-3Q17

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-05

File ID: 170802 2014 CR6-008

Sampled: 08/02/17 10:00

Prepared: 08/02/17 20:14

Analyzed: 08/02/17 20:14

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0293

Sequence: 1713798

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0020	1		EPA-7196

86092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-5

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-06

File ID: 170802 2014 CR6-009

Sampled: 08/02/17 11:01

Prepared: 08/02/17 20:14

Analyzed: 08/02/17 20:14

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0293

Sequence: 1713798

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8692117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:34:03AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-15

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-07</u>	File ID: <u>170802 2014 CR6-010</u>	
Sampled: <u>08/02/17 11:50</u>	Prepared: <u>08/02/17 20:14</u>	Analyzed: <u>08/02/17 20:14</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>BIH0293</u>	Sequence: <u>1713798</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/30/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/30/2017 10:34:03AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-8

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21317</u>
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>
Matrix: <u>Water</u>	Laboratory ID: <u>1721317-08</u>
Sampled: <u>08/02/17 13:01</u>	File ID: <u>170802 2014 CR6-018</u>
Solids: <u>0.00</u>	Prepared: <u>08/02/17 20:14</u>
Batch: <u>BIH0295</u>	Analyzed: <u>08/02/17 20:20</u>
Sequence: <u>1713798</u>	Initial/Final: <u>20 ml / 20 ml</u>
	Preparation: <u>No Prep</u>
	Calibration: <u>UNASSIGNED</u>
	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00083	1	J	EPA-7196

7/09/21/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/30/2017 10:34:03AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

EB-8-080217

Laboratory: BC Laboratories

SDG: 17-21317

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721317-09

File ID: 170802 2014 CR6-022

Sampled: 08/02/17 13:15

Prepared: 08/02/17 20:14

Analyzed: 08/02/17 20:21

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0295

Sequence:

1713798

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

1709217

LDC #: 39385C6

VALIDATION COMPLETENESS WORKSHEET

Date: 9/18/17

SDG #: 17-21317

Level III

Page: 1 of 2

Laboratory: BC Laboratories, Inc.

Reviewer: JS2nd Reviewer: JK

METHOD: (Analyte) Chloride, Nitrate-N, Sulfate (EPA Method 300.0), Nitrite-N (EPA Method 353.2), Hexavalent Chromium (EPA SW846 Method 7196), Orthophosphate-P (EPA Method 365.1), Perchlorate (EPA Method 314.0).

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A/A	
II.	Initial calibration	A	
III.	Calibration verification	A	
IV.	Laboratory Blanks	SW	
V.	Field blanks	NP	EB=8
VI.	Matrix Spike/Matrix Spike Duplicates	A	
VII.	Duplicate sample analysis	A	
VIII.	Laboratory control samples	A	LCS
IX.	Field duplicates	SW	(1,2) (3,4)
X.	Sample result verification	N	
XI.	Overall assessment of data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinse
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	MW-13	1721317-02	Water	08/02/17
2	DUP-6-3Q17	1721317-03	Water	08/02/17
3	MW-6	1721317-04	Water	08/02/17
4	DUP-7-3Q17	1721317-05	Water	08/02/17
5	MW-5	1721317-06	Water	08/02/17
6	MW-15	1721317-07	Water	08/02/17
7	MW-8	1721317-08	Water	08/02/17
8	EB-8-080217	1721317-09	Water	08/02/17
9	MW-5MS †	1721317-06MS	Water	08/02/17
10	MW-5MSD	1721317-06MSD	Water	08/02/17
11	MW-5DUP	1721317-06DUP	Water	08/02/17
12	MW-15MS †	1721317-07MS	Water	08/02/17
13	MW-15MSD	1721317-07MSD	Water	08/02/17
14	MW-15DUP	1721317-07DUP	Water	08/02/17
15	MW-8MS †	1721317-08MS	Water	08/02/17
16	MW-8MSD	1721317-08MSD	Water	08/02/17
17	MW-8DUP	1721317-08DUP	Water	08/02/17

LDC #: 39385C6

VALIDATION COMPLETENESS WORKSHEET

Date: 9/18/17

SDG #: 17-21317

Level III

Page: 2 of 2

Laboratory: BC Laboratories, Inc.

Reviewer: SB

2nd Reviewer: [Signature]

METHOD: (Analyte) Chloride, Nitrate-N, Sulfate (EPA Method 300.0), Nitrite-N (EPA Method 353.2), Hexavalent Chromium (EPA SW846 Method 7196), Orthophosphate-P (EPA Method 365.1), Perchlorate (EPA Method 314.0).

	Client ID	Lab ID	Matrix	Date
18				
19				
20				
21				
22				

Notes: _____

LDC #: 39385C6

VALIDATION FINDINGS WORKSHEET

Blanks

Page: 1 of 1

Reviewer: JB
 2nd Reviewer: AC

METHOD: Inorganics, Method See Cover

Conc. units: mg/L

Associated Samples: 1, 7

Analyte	Blank ID	Blank ID	Blank Action Limit													
	PB	ICB/CCB (mg/L)		1	7											
Nitrite as N	0.012280	0.012586	0.06293	0.015	0.014											

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:
 All contaminants within five times the method blank concentration were qualified as not detected, "U".

VALIDATION FINDINGS WORKSHEET
Field Duplicates

Inorganics, Method See Cover

Analyte	Concentration		RPD	
	1	2		
Perchlorate (ug/L)	220	230	4	
Hexavalent Chromium (mg/L)	0.0025	0.0027	8	

Analyte	Concentration		RPD	
	3	4		
Perchlorate (ug/L)	2.7	2.3	16	
Hexavalent Chromium (mg/L)	0.00070U	0.0020	96	

V:\FIELD DUPLICATES\Field Duplicates\FD_inorganic\2017\39385C6.wpd



LABORATORY DATA CONSULTANTS, INC.

2701 Loker Ave. West, Suite 220, Carlsbad, CA 92010 Bus: 760-827-1100 Fax: 760-827-1099

Tidewater, Inc.
3761 Attucks Drive
Powell, OH 43065
ATTN: Mr. David Conner

September 22, 2017

SUBJECT: NASA JPL, 3Q2017, Data Validation

Dear Mr. Conner,

Enclosed are the final validation reports for the fractions listed below. These SDGs were received on September 1, 2017. Attachment 1 is a summary of the samples that were reviewed for each analysis.

LDC Project #39394:

SDG #

Fraction

17-21091, 17-21528

Volatiles, Chromium, Wet Chemistry

The data validation was performed under Level III & IV guidelines. The analyses were validated using the following documents, as applicable to each method:

- USEPA, Contract Laboratory Program National Functional Guidelines for Superfund Organic Methods Data Review, June 2008
- USEPA, Contract Laboratory Program National Functional Guidelines for Inorganic Superfund Data Review, January 2010
- EPA SW 846, Third Edition, Test Methods for Evaluating Solid Waste, update 1, July 1992; update IIA, August 1993; update II, September 1994; update IIB, January 1995; update III, December 1996; update IIIA, April 1998; IIIB, November 2004; update IV, February 2007, update V, July 2014

Please feel free to contact us if you have any questions.

Sincerely,

Pei Geng
Project Manager/Senior Chemist

90/10 (client select)

LDC #39394 (Tidewater- Powell, OH / NASA JPL, 3Q2017)

LDC	SDG#	DATE REC'D	(3) DATE DUE	VOA (524.2)		Cr (200.8)		Cl,SO ₄ NO ₃ -N (300.0)		NO ₂ -N (353.2)		O-PO ₄ -P (365.1)		ClO ₄ (314.0)		Cr(VI) (7196)																					
				W	S	W	S	W	S	W	S	W	S	W	S	W	S	W	S	W	S	W	S	W	S	W	S	W	S	W	S	W	S	W	S		
Matrix: Water/Soil																																					
A	17-21091	09/01/17	09/25/17	12	0	9	0	1	0	1	0	1	0	11	0	9	0																				
A	17-21091	09/01/17	09/25/17	1	0	0	0	0	0	0	0	0	0	1	0	0	0																				
B	17-21528	09/01/17	09/25/17	5	0	4	0	2	0	2	0	2	0	4	0	4	0																				
Total	T/PG			18	0	13	0	3	0	3	0	3	0	16	0	13	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0	69	

Shaded cells indicate Level IV validation (all other cells are Level III validation). These sample counts do not include MS/MSD, and DUPs

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017

LDC Report Date: September 20, 2017

Parameters: Volatiles

Validation Level: Level III & IV

Laboratory: BC Laboratories, Inc.

Sample Delivery Group (SDG): 17-21091

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
TB-7-080117	1721091-01	Water	08/01/17
MW-12-5	1721091-02	Water	08/01/17
MW-12-4**	1721091-03**	Water	08/01/17
MW-12-3	1721091-04	Water	08/01/17
DUP-4-3Q17	1721091-05	Water	08/01/17
MW-12-2	1721091-06	Water	08/01/17
MW-12-1	1721091-07	Water	08/01/17
MW-11-4	1721091-08	Water	08/01/17
MW-11-3	1721091-09	Water	08/01/17
MW-11-2	1721091-10	Water	08/01/17
DUP-5-3Q17	1721091-11	Water	08/01/17
MW-11-1	1721091-12	Water	08/01/17
EB-7-080117	1721091-13	Water	08/01/17
MW-12-5MS	1721091-02MS	Water	08/01/17
MW-12-5MSD	1721091-02MSD	Water	08/01/17

**Indicates sample underwent Level IV review

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Superfund Organic Methods Data Review (June 2008). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following method:

Volatile Organic Compounds (VOCs) by Environmental Protection Agency (EPA) Method 524.2

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results. Samples appended with a double asterisk on the cover page were subjected to Level IV evaluation, which is comprised of the QC summary forms as well as the raw data, to confirm sample quantitation and identification.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition and cooler temperatures upon receipt met validation criteria.

All technical holding time requirements were met.

II. GC/MS Instrument Performance Check

Instrument performance was checked at 12 hour intervals.

All ion abundance requirements were met.

III. Initial Calibration and Initial Calibration Verification

An initial calibration was performed as required by the method.

For compounds where average relative response factors (RRFs) were utilized, the percent relative standard deviations (%RSD) were less than or equal to 20.0%.

In the case where the laboratory used a calibration curve to evaluate the compounds, all coefficients of determination (r^2) were greater than or equal to 0.990.

The percent differences (%D) of the initial calibration verification (ICV) standard were less than or equal to 30.0% for all compounds.

IV. Continuing Calibration

Continuing calibration was performed at the required frequencies.

The percent differences (%D) were less than or equal to 30.0% for all compounds with the following exceptions:

Date	Compound	%D	Associated Samples	Flag	A or P
08/04/17	Methyl iodide Pentachloroethane	51.8 (≤ 30) 33.9 (≤ 30)	All samples in SDG 17-21091	UJ (all non-detects) UJ (all non-detects)	P

V. Laboratory Blanks

Laboratory blanks were analyzed as required by the method. No contaminants were found in the laboratory blanks.

VI. Field Blanks

Sample TB-7-080117 was identified as a trip blank. No contaminants were found.

Sample EB-7-080117 was identified as an equipment blank. No contaminants were found.

VII. Surrogates

Surrogates were added to all samples as required by the method. All surrogate recoveries (%R) were within QC limits.

VIII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

IX. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the method. Percent recoveries (%R) were within QC limits.

X. Field Duplicates

Samples MW-12-3 and DUP-4-3Q17 and samples MW-11-2 and DUP-5-3Q17 were identified as field duplicates. No results were detected in any of the samples with the following exceptions:

Analyte	Concentration (ug/L)		RPD
	MW-12-3	DUP-4-3Q17	
Carbon tetrachloride	0.40	0.32	22
Chloroform	0.78	0.63	21

XI. Internal Standards

All internal standard areas and retention times were within QC limits.

XII. Compound Quantitation

All compound quantitations met validation criteria for samples which underwent Level IV validation. Raw data were not reviewed for Level III validation.

XIII. Target Compound Identifications

All target compound identifications met validation criteria for samples which underwent Level IV validation. Raw data were not reviewed for Level III validation.

XIV. System Performance

The system performance was acceptable for samples which underwent Level IV validation. Raw data were not reviewed for Level III validation.

XV. Overall Assessment of Data

The analysis was conducted within all specifications of the method. No results were rejected in this SDG.

Due to continuing calibration %D, data were qualified as estimated in thirteen samples.

The quality control criteria reviewed, other than those discussed above, were met and are considered acceptable. Sample results that were found to be estimated (J) are usable for limited purposes only. Based upon the data validation all other results are considered valid and usable for all purposes.

NASA JPL, 3Q2017

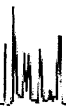
Volatiles - Data Qualification Summary - SDG 17-21091

Sample	Compound	Flag	A or P	Reason
TB-7-080117 MW-12-5 MW-12-4** MW-12-3 DUP-4-3Q17 MW-12-2 MW-12-1 MW-11-4 MW-11-3 MW-11-2 DUP-5-3Q17 MW-11-1 EB-7-080117	Methyl iodide Pentachloroethane	UJ (all non-detects) UJ (all non-detects)	P	Continuing calibration (%D)

NASA JPL, 3Q2017

Volatiles - Laboratory Blank Data Qualification Summary - SDG 17-21091

No Sample Data Qualified in this SDG



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

TB-7-080117

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-01</u>	File ID:	<u>04AUG14.D</u>		
Sampled:	<u>08/01/17 07:00</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 14:17</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>B[H0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/22/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

TB-7-080117

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-01 File ID: 04AUG14.D
Sampled: 08/01/17 07:00 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 14:17
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	8.3600	83.6	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.2000	92.0	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	11.220	112	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	223610	6.58	230753	6.57	
Chlorobenzene-d5 (IS)	76251	9.61	81486	9.61	
1,4-Difluorobenzene (IS)	312335	7.38	338008	7.38	

* Values outside of QC limits

8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET

EPA-524.2

TB-7-080117

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-01</u>	File ID:	<u>04AUG14.D</u>		
Sampled:	<u>08/01/17 07:00</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 14:17</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BJH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

17572117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-12-5

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-02 File ID: 04AUG07.D
Sampled: 08/01/17 07:30 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 11:36
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.28	J
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-12-5

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-02</u>	File ID:	<u>04AUG07.D</u>		
Sampled:	<u>08/01/17 07:30</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 11:36</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BJH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

11092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-12-5

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-02</u>	File ID:	<u>04AUG07.D</u>		
Sampled:	<u>08/01/17 07:30</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 11:36</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>U5</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>U5</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	9.6400	96.4	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.3200	93.2	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	8.7700	87.7	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	221283	6.57	230753	6.57	
Chlorobenzene-d5 (IS)	79776	9.62	81486	9.61	
1,4-Difluorobenzene (IS)	303623	7.38	338008	7.38	

* Values outside of QC limits

1092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-12-5

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-02</u>	File ID:	<u>04AUG07.D</u>		
Sampled:	<u>08/01/17 07:30</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 11:36</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BJH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8692117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-12-4

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-03</u>	File ID: <u>04AUG15.D</u>	
Sampled: <u>08/01/17 08:00</u>	Prepared: <u>08/04/17 07:00</u>	Analyzed: <u>08/04/17 14:40</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BJH0384</u>	Sequence: <u>1713791</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.30	J
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.41	J
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/21/17

Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-12-4

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-03</u>	File ID: <u>04AUG15.D</u>	
Sampled: <u>08/01/17 08:00</u>	Prepared: <u>08/04/17 07:00</u>	Analyzed: <u>08/04/17 14:40</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BJH0384</u>	Sequence: <u>1713791</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>U5</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>U5</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	9.8300	98.3	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.8100	98.1	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.1600	91.6	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	215534	6.57	230753	6.57	
Chlorobenzene-d5 (IS)	78546	9.61	81486	9.61	
1,4-Difluorobenzene (IS)	329340	7.38	338008	7.38	

* Values outside of QC limits

82092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET

EPA-524.2

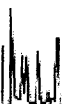
MW-12-4

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-03</u>	File ID: <u>04AUG15.D</u>	
Sampled: <u>08/01/17 08:00</u>	Prepared: <u>08/04/17 07:00</u>	Analyzed: <u>08/04/17 14:40</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BH0384</u>	Sequence: <u>1713791</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-12-3

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-04 File ID: 04AUG16.D
Sampled: 08/01/17 08:15 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 15:03
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.40	J
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.78	
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

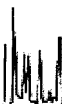
ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-12-3

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-04 File ID: 04AUG16.D
Sampled: 08/01/17 08:15 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 15:03
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

209217



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-12-3

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-04 File ID: 04AUG16.D
Sampled: 08/01/17 08:15 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 15:03
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

Table with 5 columns: CAS NO., COMPOUND, DILUTION, CONC. (ug/L), Q. Lists various organic compounds and their concentrations.

Table with 6 columns: SYSTEM MONITORING COMPOUND, ADDED (ug/L), CONC (ug/L), % REC, QC LIMITS, Q. Shows surrogate compounds and their recovery percentages.

Table with 6 columns: INTERNAL STANDARD, AREA, RT, REF AREA, REF RT, Q. Lists internal standards used for calibration.

* Values outside of QC limits

Signature



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET

EPA-524.2

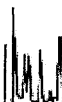
MW-12-3

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-04</u>	File ID:	<u>04AUG16.D</u>		
Sampled:	<u>08/01/17 08:15</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 15:03</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

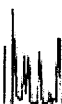
EPA-524.2

DUP-4-3Q17

Laboratory: BC Laboratories SDG: 17-21091
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1721091-05 File ID: 04AUG17.D
 Sampled: 08/01/17 08:25 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 15:26
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: B/H0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.32	J
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.63	
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/31/17


 Tidewater Inc.
 3761 Attucks Drive
 Powell, OH 43065

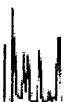
 Reported: 8/31/2017 11:08:10AM
 Project: JPL- GW Monitoring Wells
 Project Number: 3Q17
 Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2
DUP-4-3Q17

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-05</u>
		File ID:	<u>04AUG17.D</u>
Sampled:	<u>08/01/17 08:25</u>	Prepared:	<u>08/04/17 07:00</u>
		Analyzed:	<u>08/04/17 15:26</u>
Solids:		Preparation:	<u>EPA 5030 Water MS</u>
		Initial/Final:	<u>25 ml / 25 ml</u>
Batch:	<u>BIH0384</u>	Sequence:	<u>1713791</u>
		Calibration:	<u>1707017</u>
		Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

DUP-4-3Q17

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-05</u>	File ID:	<u>04AUG17.D</u>		
Sampled:	<u>08/01/17 08:25</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 15:26</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>V5</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>V5</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	9.8900	98.9	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.7300	97.3	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	8.6200	86.2	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	210733	6.57	230753	6.57	
Chlorobenzene-d5 (IS)	81341	9.61	81486	9.61	
1,4-Difluorobenzene (IS)	317100	7.38	338008	7.38	

* Values outside of QC limits

Step 17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET

EPA-524.2

DUP-4-3Q17

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-05</u>	File ID:	<u>04AUG17.D</u>		
Sampled:	<u>08/01/17 08:25</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 15:26</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BJH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-12-2

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-06 File ID: 04AUG18.D
Sampled: 08/01/17 09:15 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 15:49
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-12-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-06</u>	File ID: <u>04AUG18.D</u>	
Sampled: <u>08/01/17 09:15</u>	Prepared: <u>08/04/17 07:00</u>	Analyzed: <u>08/04/17 15:49</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BJH0384</u>	Sequence: <u>1713791</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-12-2

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-06</u>	File ID: <u>04AUG18.D</u>	
Sampled: <u>08/01/17 09:15</u>	Prepared: <u>08/04/17 07:00</u>	Analyzed: <u>08/04/17 15:49</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BJH0384</u>	Sequence: <u>1713791</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>U5</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>U5</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.310	103	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.9000	99.0	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.7700	97.7	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	214627	6.57	230753	6.57	
Chlorobenzene-d5 (IS)	85142	9.61	81486	9.61	
1,4-Difluorobenzene (IS)	318108	7.38	338008	7.38	

* Values outside of QC limits

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Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

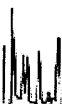
EPA-524.2

MW-12-1

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-07 File ID: 04AUG19.D
Sampled: 08/01/17 09:45 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 16:12
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B/H0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-12-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-07</u>	File ID: <u>04AUG19.D</u>	
Sampled: <u>08/01/17 09:45</u>	Prepared: <u>08/04/17 07:00</u>	Analyzed: <u>08/04/17 16:12</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BJH0384</u>	Sequence: <u>1713791</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-12-1

Laboratory: BC Laboratories SDG: 17-21091
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1721091-07 File ID: 04AUG19.D
 Sampled: 08/01/17 09:45 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 16:12
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

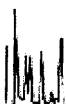
CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.380	104	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.8200	98.2	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.8200	98.2	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	209239	6.57	230753	6.57	
Chlorobenzene-d5 (IS)	83261	9.62	81486	9.61	
1,4-Difluorobenzene (IS)	311271	7.38	338008	7.38	

* Values outside of QC limits

8/30/2017



Tidewater Inc.
 3761 Attucks Drive
 Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
 Project: JPL- GW Monitoring Wells
 Project Number: 3Q17
 Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-4

Laboratory: BC Laboratories SDG: 17-21091
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1721091-08 File ID: 04AUG20.D
 Sampled: 08/01/17 11:00 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 16:35
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-11-4

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-08 File ID: 04AUG20.D
Sampled: 08/01/17 11:00 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 16:35
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

509217



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-4

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-08 File ID: 04AUG20.D
Sampled: 08/01/17 11:00 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 16:35
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>US</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>US</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.410	104	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.7400	97.4	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.6000	96.0	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	205701	6.58	230753	6.57	
Chlorobenzene-d5 (IS)	85483	9.62	81486	9.61	
1,4-Difluorobenzene (IS)	316984	7.38	338008	7.38	

* Values outside of QC limits

5607217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-4

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-08</u>	File ID:	<u>04AUG20.D</u>		
Sampled:	<u>08/01/17 11:00</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 16:35</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>B[H0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

Signature



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-11-3

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-09 File ID: 04AUG21.D
Sampled: 08/01/17 11:30 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 16:58
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-3

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-09 File ID: 04AUG21.D
Sampled: 08/01/17 11:30 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 16:58
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.13	J
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-3

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-09 File ID: 04AUG21.D
Sampled: 08/01/17 11:30 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 16:58
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	10.210	102	75 - 125	
Toluene-d8 (Surrogate)	10.000	8.7100	87.1	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.6400	96.4	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	210850	6.57	230753	6.57	
Chlorobenzene-d5 (IS)	76849	9.62	81486	9.61	
1,4-Difluorobenzene (IS)	317066	7.38	338008	7.38	

* Values outside of QC limits

82092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-3

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-09</u>	File ID: <u>04AUG21.D</u>	
Sampled: <u>08/01/17 11:30</u>	Prepared: <u>08/04/17 07:00</u>	Analyzed: <u>08/04/17 16:58</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BIH0384</u>	Sequence: <u>1713791</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

20592117



Tidewater Inc.
 3761 Attucks Drive
 Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
 Project: JPL- GW Monitoring Wells
 Project Number: 3Q17
 Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-11-2

Laboratory: BC Laboratories SDG: 17-21091
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1721091-10 File ID: 04AUG22.D
 Sampled: 08/01/17 11:50 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 17:21
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: BIH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-2

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-10 File ID: 04AUG22.D
Sampled: 08/01/17 11:50 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 17:21
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-2

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-10 File ID: 04AUG22.D
Sampled: 08/01/17 11:50 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 17:21
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
75-15-0	Carbon disulfide	1	0.48	U
110-57-6	trans-1,4-Dichloro-2-butene	1	1.8	U
60-29-7	Diethyl ether	1	0.33	U
97-63-2	Ethyl methacrylate	1	1.3	U
637-92-3	Ethyl t-butyl ether	1	0.32	U
67-72-1	Hexachloroethane	1	0.11	U
591-78-6	2-Hexanone	1	5.0	U
126-98-7	Methacrylonitrile	1	2.3	U
78-93-3	Methyl ethyl ketone	1	3.3	U
74-88-4	Methyl iodide	1	1.1	U <i>UJ</i>
108-10-1	Methyl isobutyl ketone	1	2.4	U
80-62-6	Methyl methacrylate	1	1.2	U
76-01-7	Pentachloroethane	1	0.63	U <i>UJ</i>
107-12-0	Propionitrile	1	6.2	U
109-99-9	Tetrahydrofuran	1	5.2	U
179601-23-1	p- & m-Xylenes	1	0.34	U
95-47-6	o-Xylene	1	0.13	U

SYSTEM MONITORING COMPOUND	ADDED (ug/L)	CONC (ug/L)	% REC	QC LIMITS	Q
1,2-Dichloroethane-d4 (Surrogate)	10.000	9.6900	96.9	75 - 125	
Toluene-d8 (Surrogate)	10.000	9.0900	90.9	80 - 120	
4-Bromofluorobenzene (Surrogate)	10.000	9.2600	92.6	80 - 120	

INTERNAL STANDARD	AREA	RT	REF AREA	REF RT	Q
Pentafluorobenzene (IS)	205299	6.58	230753	6.57	
Chlorobenzene-d5 (IS)	73221	9.61	81486	9.61	
1,4-Difluorobenzene (IS)	296757	7.38	338008	7.38	

* Values outside of QC limits

8059217



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

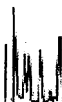
MW-11-2

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-10 File ID: 04AUG22.D
Sampled: 08/01/17 11:50 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 17:21
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

86092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

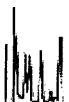
EPA-524.2

DUP-5-3Q17

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-11 File ID: 04AUG23.D
Sampled: 08/01/17 12:00 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 17:44
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

82092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

DUP-5-3Q17

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-11 File ID: 04AUG23.D
Sampled: 08/01/17 12:00 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 17:44
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BJH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

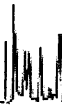
Table with 5 columns: CAS NO., COMPOUND, DILUTION, CONC. (ug/L), Q. Lists various organic compounds and their concentrations.

Table with 6 columns: SYSTEM MONITORING COMPOUND, ADDED (ug/L), CONC (ug/L), % REC, QC LIMITS, Q. Shows surrogate compounds and their recovery percentages.

Table with 6 columns: INTERNAL STANDARD, AREA, RT, REF AREA, REF RT, Q. Lists internal standards used for calibration.

* Values outside of QC limits

Signature: S092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

DUP-5-3Q17

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-11 File ID: 04AUG23.D
Sampled: 08/01/17 12:00 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 17:44
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B\H0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-11-1

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-12</u>	File ID:	<u>04AUG24.D</u>		
Sampled:	<u>08/01/17 12:50</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 18:07</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

Signature



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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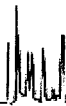
ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-1

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-12</u>	File ID:	<u>04AUG24.D</u>		
Sampled:	<u>08/01/17 12:50</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 18:07</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BJH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

82092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-1

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-12 File ID: 04AUG24.D
Sampled: 08/01/17 12:50 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 18:07
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

Table with 5 columns: CAS NO., COMPOUND, DILUTION, CONC. (ug/L), Q. Lists various organic compounds and their concentrations.

Table with 6 columns: SYSTEM MONITORING COMPOUND, ADDED (ug/L), CONC (ug/L), % REC, QC LIMITS, Q. Shows surrogate compounds and their recovery percentages.

Table with 6 columns: INTERNAL STANDARD, AREA, RT, REF AREA, REF RT, Q. Lists internal standards used for calibration.

* Values outside of QC limits

Signature



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-11-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-12</u>	File ID: <u>04AUG24.D</u>	
Sampled: <u>08/01/17 12:50</u>	Prepared: <u>08/04/17 07:00</u>	Analyzed: <u>08/04/17 18:07</u>	
Solids:	Preparation: <u>EPA 5030 Water MS</u>	Initial/Final: <u>25 ml / 25 ml</u>	
Batch: <u>BJH0384</u>	Sequence: <u>1713791</u>	Calibration: <u>1707017</u>	Instrument: <u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET

EPA-524.2

EB-7-080117

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-13 File ID: 04AUG25.D
Sampled: 08/01/17 13:00 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 18:30
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: BIH0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:08:10AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-7-080117

Laboratory: BC Laboratories SDG: 17-21091
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721091-13 File ID: 04AUG25.D
Sampled: 08/01/17 13:00 Prepared: 08/04/17 07:00 Analyzed: 08/04/17 18:30
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0384 Sequence: 1713791 Calibration: 1707017 Instrument: MS-V5

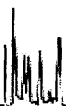
Table with 5 columns: CAS NO., COMPOUND, DILUTION, CONC. (ug/L), Q. Lists various organic compounds and their concentrations.

Table with 6 columns: SYSTEM MONITORING COMPOUND, ADDED (ug/L), CONC (ug/L), % REC, QC LIMITS, Q. Lists surrogate compounds and their recovery percentages.

Table with 6 columns: INTERNAL STANDARD, AREA, RT, REF AREA, REF RT, Q. Lists internal standards and their retention times.

* Values outside of QC limits

Handwritten signature/initials



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:08:10AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET

EPA-524.2

EB-7-080117

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21091</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721091-13</u>	File ID:	<u>04AUG25.D</u>		
Sampled:	<u>08/01/17 13:00</u>	Prepared:	<u>08/04/17 07:00</u>	Analyzed:	<u>08/04/17 18:30</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BJH0384</u>	Sequence:	<u>1713791</u>	Calibration:	<u>1707017</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

Signature

LDC #: 39394A1

VALIDATION COMPLETENESS WORKSHEET

SDG #: 17-21091

Level III/IV

Laboratory: BC Laboratories, Inc.

Date: 9/12/17

Page: 1 of 2

Reviewer: 9

2nd Reviewer: R

METHOD: GC/MS Volatiles (EPA Method 524.2)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A	
II.	GC/MS Instrument performance check	A	
III.	Initial calibration/ICV	AA	RSD ≤ 20% γ^2 1CV ≤ 3%
IV.	Continuing calibration	M	CCV ≤ 3%
V.	Laboratory Blanks	A	
VI.	Field blanks	ND	TB = 1. 2/3 = 1/3
VII.	Surrogate spikes	A	
VIII.	Matrix spike/Matrix spike duplicates	A	
IX.	Laboratory control samples	A	LCS
X.	Field duplicates	M	D = 4+5. 10+11
XI.	Internal standards	A	
XII.	Compound quantitation RL/LOQ/LODs	A	Not reviewed for Level III validation.
XIII.	Target compound identification	A	Not reviewed for Level III validation.
XIV.	System performance	A	Not reviewed for Level III validation.
XV.	Overall assessment of data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

** Indicates sample was underwent Level IV review

	Client ID	Lab ID	Matrix	Date
1	TB-7-080117	1721091-01	Water	08/01/17
2	MW-12-5	1721091-02	Water	08/01/17
3	MW-12-4**	1721091-03**	Water	08/01/17
4	MW-12-3	1721091-04	Water	08/01/17
5	DUP-4-3Q17	1721091-05	Water	08/01/17
6	MW-12-2	1721091-06	Water	08/01/17
7	MW-12-1	1721091-07	Water	08/01/17
8	MW-11-4	1721091-08	Water	08/01/17
9	MW-11-3	1721091-09	Water	08/01/17
10	MW-11-2	1721091-10	Water	08/01/17
11	DUP-5-3Q17	1721091-11	Water	08/01/17
12	MW-11-1	1721091-12	Water	08/01/17
13	EB-7-080117	1721091-13	Water	08/01/17

LDC #: 39394A1

VALIDATION COMPLETENESS WORKSHEET

SDG #: 17-21091

Level III/IV

Laboratory: BC Laboratories, Inc.

Date: 9/12/17

Page: 2 of 2

Reviewer: J

2nd Reviewer: H

METHOD: GC/MS Volatiles (EPA Method 524.2)

	Client ID	Lab ID	Matrix	Date
14	MW-12-5MS	1721091-02MS	Water	08/01/17
15	MW-12-5MSD	1721091-02MSD	Water	08/01/17
16				
17				
18				
19				
20				

Notes:

Method: Volatiles (EPA Method 524.2)

Validation Area	Yes	No	NA	Findings/Comments
I. Technical holding times				
Were all technical holding times met?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was cooler temperature criteria met?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
II. GC/MS Instrument performance check				
Was a tune check performed prior to establishing and/or re-establishing an initial calibration?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the BFB performance results reviewed and found to be within the specified criteria?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
III. Initial calibration				
Did the laboratory perform at least 5 point calibration prior to sample analysis?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all percent relative standard deviations (%RSD) < 20%?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
IIIa. Initial Calibration Verification calibration				
Was an initial calibration verification standard analyzed after each initial calibration for each instrument?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all percent differences (%D) < 30%?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
IV. Continuing calibration				
Was a continuing calibration standard analyzed at the beginning of each analysis batch?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were all percent differences (%D) of continuing calibration < 30%?	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
V. Laboratory Blanks				
Was a laboratory blank associated with every sample in this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was a laboratory blank analyzed with each analysis batch?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Was there contamination in the laboratory blanks? If yes, please see the Blanks validation completeness worksheet.	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
VI. Field blanks				
Field blanks were identified in this SDG.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Target compounds were detected in the field blanks.	<input type="checkbox"/>	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
VII. Surrogate spikes				
Were all surrogate %R within the QC limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
If the percent recovery (%R) for one or more surrogates was out of QC limits, was a reanalysis performed to confirm samples with %R outside of criteria?	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
VIII. Matrix spike/Matrix spike duplicates				
Was a matrix spike (MS) and matrix spike duplicate (MSD) analyzed for this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the QC limits?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
IX. Laboratory control samples				
Was an LCS analyzed for this SDG?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	

VALIDATION FINDINGS CHECKLIST

Validation Area	Yes	No	NA	Findings/Comments
Was an LCS analyzed per analytical batch?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were the LCS percent recoveries (%R) within 70-130%?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
X. Field duplicates				
Field duplicate pairs were identified in this SDG.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Target compounds were detected in the field duplicates.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
XI. Internal standards				
Were internal standard area counts within +/-30% of the area of the most recent continuing calibration standard and +/-50% of the average peak area in the initial calibration?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were retention times within +/-30 seconds of the associated calibration standard?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
XII. Compound quantitation/CRQLs				
Were the correct internal standard (IS), quantitation ion and relative response factor (RRF) or regression equations used to quantitate the compound?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were compound quantitation and CRQLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
XIII. Target compound identification				
Were relative retention times (RRT's) within + 0.06 RRT units of the standard?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Did compound spectra meet specified EPA "Functional Guidelines" criteria?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
Were chromatogram peaks verified and accounted for?	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
XIV. System performance				
System performance was found to be acceptable.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	
XV. Overall assessment of data				
Overall assessment of data was found to be acceptable.	<input checked="" type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	

TARGET COMPOUND WORKSHEET

METHOD: VOA

A. Chloromethane	AA. Tetrachloroethene	AAA. 1,3,5-Trimethylbenzene	AAAA. Ethyl tert-butyl ether	A1. 1,3-Butadiene
B. Bromomethane	BB. 1,1,2,2-Tetrachloroethane	BBB. 4-Chlorotoluene	BBBB. tert-Amyl methyl ether	B1. Hexane
C. Vinyl chloride	CC. Toluene	CCC. tert-Butylbenzene	CCCC. 1-Chlorohexane	C1. Heptane
D. Chloroethane	DD. Chlorobenzene	DDD. 1,2,4-Trimethylbenzene	DDDD. Isopropyl alcohol	D1. Propylene
E. Methylene chloride	EE. Ethylbenzene	EEE. sec-Butylbenzene	EEEE. Acetonitrile	E1. Freon 11
F. Acetone	FF. Styrene	FFF. 1,3-Dichlorobenzene	FFFF. Acrolein	F1. Freon 12
G. Carbon disulfide	GG. Xylenes, total	GGG. p-Isopropyltoluene	GGGG. Acrylonitrile	G1. Freon 113
H. 1,1-Dichloroethene	HH. Vinyl acetate	HHH. 1,4-Dichlorobenzene	HHHH. 1,4-Dioxane	H1. Freon 114
I. 1,1-Dichloroethane	II. 2-Chloroethylvinyl ether	III. n-Butylbenzene	IIII. Isobutyl alcohol	I1. 2-Nitropropane
J. 1,2-Dichloroethene, total	JJ. Dichlorodifluoromethane	JJJ. 1,2-Dichlorobenzene	JJJJ. Methacrylonitrile	J1. Dimethyl disulfide
K. Chloroform	KK. Trichlorofluoromethane	KKK. 1,2,4-Trichlorobenzene	KKKK. Propionitrile	K1. 2,3-Dimethyl pentane
L. 1,2-Dichloroethane	LL. Methyl-tert-butyl ether	LLL. Hexachlorobutadiene	LLLL. Ethyl ether	L1. 2,4-Dimethyl pentane
M. 2-Butanone	MM. 1,2-Dibromo-3-chloropropane	MMM. Naphthalene	MMMM. Benzyl chloride	M1. 3,3-Dimethyl pentane
N. 1,1,1-Trichloroethane	NN. Methyl ethyl ketone	NNN. 1,2,3-Trichlorobenzene	NNNN. Iodomethane	N1. 2-Methylpentane
O. Carbon tetrachloride	OO. 2,2-Dichloropropane	OOO. 1,3,5-Trichlorobenzene	OOOO. 1,1-Difluoroethane	O1. 3-Methylpentane
P. Bromodichloromethane	PP. Bromochloromethane	PPP. trans-1,2-Dichloroethene	PPPP. Tetrahydrofuran	P1. 3-Ethylpentane
Q. 1,2-Dichloropropane	QQ. 1,1-Dichloropropene	QQQ. cis-1,2-Dichloroethene	QQQQ. Methyl acetate	Q1. 2,2-Dimethylpentane
R. cis-1,3-Dichloropropene	RR. Dibromomethane	RRR. m,p-Xylenes	RRRR. Ethyl acetate	R1. 2,2,3-Trimethylbutane
S. Trichloroethene	SS. 1,3-Dichloropropane	SSS. o-Xylene	SSSS. Cyclohexane	S1. 2,2,4-Trimethylpentane
T. Dibromochloromethane	TT. 1,2-Dibromoethane	TTT. 1,1,2-Trichloro-1,2,2-trifluoroethane	TTTT. Methylcyclohexane	T1. 2-Methylhexane
U. 1,1,2-Trichloroethane	UU. 1,1,1,2-Tetrachloroethane	UUU. 1,2-Dichlorotetrafluoroethane	UUUU. Allyl chloride	U1. Nonanal
V. Benzene	VV. Isopropylbenzene	VVV. 4-Ethyltoluene	VVVV. Methyl methacrylate	V1. 2-Methylnaphthalene
W. trans-1,3-Dichloropropene	WW. Bromobenzene	WWW. Ethanol	WWWW. Ethyl methacrylate	W1. Methanol
X. Bromoform	XX. 1,2,3-Trichloropropane	XXX. Di-isopropyl ether	XXXX. cis-1,4-Dichloro-2-butene	X1. 1,2,3-Trimethylbenzene
Y. 4-Methyl-2-pentanone	YY. n-Propylbenzene	YYY. tert-Butanol	YYYY. trans-1,4-Dichloro-2-butene	Y1.
Z. 2-Hexanone	ZZ. 2-Chlorotoluene	ZZZ. tert-Butyl alcohol	ZZZZ. Pentachloroethane	Z1.

LDC#: 39394A1

VALIDATION FINDINGS WORKSHEET
Field Duplicates

Page: 1 of 1
Reviewer: [Signature]
2nd Reviewer: [Signature]

METHOD: GCMS VOA (EPA Method 524.2)

Compound	Concentration (ug/L)		RPD
	4	5	
O	0.40	0.32	22
K	0.78	0.63	21

V:\FIELD DUPLICATES\Field Duplicates\FD_Organics\2017\39394A1_JPL.wpd

VALIDATION FINDINGS WORKSHEET Initial Calibration Calculation Verification

METHOD: GC/MS VOA (EPA SW 846 Method 8260B)

The Relative Response Factor (RRF), average RRF, and percent relative standard deviation (%RSD) were recalculated for the compounds identified below using the following calculations:

$$RRF = (A_x)(C_{is}) / (A_{is})(C_x)$$

average RRF = sum of the RRFs/number of standards

$$\%RSD = 100 * (S/X)$$

A_x = Area of compound,

C_x = Concentration of compound,

S = Standard deviation of the RRFs

X = Mean of the RRFs

A_{is} = Area of associated internal standard

C_{is} = Concentration of internal standard

#	Standard ID	Calibration Date	Compound (Reference Internal Standard)	Reported	Recalculated	Reported	Recalculated	Reported	Recalculated
				RRF (10 std)	RRF (10 std)	Average RRF (initial)	Average RRF (initial)	%RSD	%RSD
1	ICAL (V5)	7/20/17	K (1st internal standard)	0.7797252	0.7797252	0.7571101	0.7571101	5.641054	5.641053
			S (2nd internal standard)	0.3603684	0.3603684	0.3434012	0.3434012	3.079274	3.079306
			EE (3rd internal standard)	2.063766	2.063766	1.948304	1.948304	6.711166	6.711173
			BB (4th internal standard)						
2			K (1st internal standard)						
			S (2nd internal standard)						
			AA (3rd internal standard)						
			BB (4th internal standard)						
3			K (1st internal standard)						
			S (2nd internal standard)						
			AA (3rd internal standard)						
			BB (4th internal standard)						
4			K (1st internal standard)						
			S (2nd internal standard)						
			AA (3rd internal standard)						
			BB (4th internal standard)						

Comments: Refer to Initial Calibration findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

VALIDATION FINDINGS WORKSHEET Continuing Calibration Results Verification

METHOD: GC/MS VOA (EPA Method 524.2)

The percent difference (%D) of the initial calibration average Relative Response Factors (RRFs) and the continuing calibration RRFs were recalculated for the compounds identified below using the following calculation:

% Difference = $100 * (\text{ave. RRF} - \text{RRF}) / \text{ave. RRF}$
 $\text{RRF} = (A_x)(C_{is}) / (A_{is})(C_x)$

Where: ave. RRF = initial calibration average RRF
 RRF = continuing calibration RRF
 A_x = Area of compound, A_{is} = Area of associated internal standard
 C_x = Concentration of compound, C_{is} = Concentration of internal standard

#	Standard ID	Calibration Date	Compound (Reference internal Standard)	Average RRF (initial)	Reported RRF (CC)	Recalculated RRF (CC)	Reported %D	Recalculated %D
1	04aug02	8/4/17	K (1st internal standard)	0.7571101	0.7733756	0.7733756	2.1	2.1
			S (2nd internal standard)	0.3434012	0.3649989	0.3649989	6.3	6.3
			EE (3rd internal standard)	1.948304	2.011687	2.011687	3.3	3.3
			BB (4th internal standard)					
2			K (5th internal standard)					
			S (2nd internal standard)					
			AA (3rd internal standard)					
			BB (4th internal standard)					
3			QQQ (1st internal standard)					
			S (2nd internal standard)					
			AA (3rd internal standard)					
			MMM (4th internal standard)					
4			BB (5th internal standard)					
			S (2nd internal standard)					
			AA (3rd internal standard)					
			BB (4th internal standard)					

Comments: Refer to Continuing Calibration findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

VALIDATION FINDINGS WORKSHEET
Surrogate Results Verification

METHOD: GC/MS VOA (EPA Method 524.2)

The percent recoveries (%R) of surrogates were recalculated for the compounds identified below using the following calculation:

% Recovery: $SF/SS * 100$

Where: SF = Surrogate Found
 SS = Surrogate Spiked

Sample ID: 3

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8	10.000	9.81	98.1	98.1	0
Bromofluorobenzene	↓	9.16	91.6	91.6	↓
1,2-Dichlorobenzene-d4	↓	98.3	98.3	98.3	↓
Dibromofluoromethane					

Sample ID: _____

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					
Dibromofluoromethane					

Sample ID: _____

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					
Dibromofluoromethane					

Sample ID: _____

	Surrogate Spiked	Surrogate Found	Percent Recovery	Percent Recovery	Percent Difference
			Reported	Recalculated	
Toluene-d8					
Bromofluorobenzene					
1,2-Dichlorobenzene-d4					
Dibromofluoromethane					

VALIDATION FINDINGS WORKSHEET
Matrix Spike/Matrix Spike Duplicates Results Verification

METHOD: GC/MS VOA (EPA Method 524.2)

The percent recoveries (%R) and Relative Percent Difference (RPD) of the matrix spike and matrix spike duplicate were recalculated for the compounds identified below using the following calculation:

% Recovery = 100 * (SSC - SC)/SA

Where: SSC = Spiked sample concentration
 SA = Spike added

SC = Sample concentration

RPD = | MSC - MSC | * 2 / (MSC + MSDC)

MSC = Matrix spike concentration

MSDC = Matrix spike duplicate concentration

MS/MSD sample: 14/15

Compound	Spike Added		Sample Concentration	Spiked Sample Concentration		Matrix Spike		Matrix Spike Duplicate		MS/MSD	
	MS	MSD		MS	MSD	Percent Recovery		Percent Recovery		RPD	
						Reported	Recalc	Reported	Recalc	Reported	Recalculated
1,1-Dichloroethene	25.000	25.000	ND	29.760	26.750	119	119	105	105	12.2	12.2
Trichloroethene	↓	↓	↓	27.320	26.210	109	109	105	105	4.15	4.15
Benzene	↓	↓	↓	22.850	24.750	91.4	91.4	97.0	97.0	5.94	5.94
Toluene	↓	↓	↓	27.150	25.800	109	109	103	103	5.10	5.10
Chlorobenzene	↓	↓	↓	22.550	22.730	90.2	90.2	90.9	90.9	0.795	0.795

Comments: Refer to Matrix Spike/Matrix Spike Duplicates findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

VALIDATION FINDINGS WORKSHEET
Laboratory Control Sample Results Verification

METHOD: GC/MS VOA (EPA Method 524.2)

The percent recoveries (%R) and Relative Percent Difference (RPD) of the laboratory control sample and laboratory control sample duplicate (if applicable) were recalculated for the compounds identified below using the following calculation:

% Recovery = 100 * SSC/SA

Where: SSC = Spiked sample concentration
 SA = Spike added

RPD = |LCS - LCSDC| * 2 / (LCS + LCSDC)

LCS = Laboratory control sample concentration LCSDC = Laboratory control sample duplicate concentration

LCS ID: B1H0387-351

Compound	Spike Added (<u>1000</u>)		Spiked Sample Concentration (<u>1000</u>)		LCS		LCSD		LCS/LCSD	
	LCS	LCSD	LCS	LCSD	Percent Recovery		Percent Recovery		RPD	
					Reported	Recalc.	Reported	Recalc.	Reported	Recalculated
1,1-Dichloroethene	<u>25.00</u>	<u>NA</u>	<u>26.790</u>	<u>NA</u>	<u>107</u>	<u>107</u>				
Trichloroethene	↓	↓	<u>25.710</u>	↓	<u>103</u>	<u>103</u>				
Benzene	↓	↓	<u>24.370</u>	↓	<u>97.5</u>	<u>97.5</u>				
Toluene	↓	↓	<u>23.760</u>	↓	<u>95.0</u>	<u>95.0</u>				
Chlorobenzene	↓	↓	<u>23.320</u>	↓	<u>93.3</u>	<u>93.3</u>				

Comments: Refer to Laboratory Control Sample findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017

LDC Report Date: September 21, 2017

Parameters: Chromium

Validation Level: Level III

Laboratory: BC Laboratories, Inc.

Sample Delivery Group (SDG): 17-21091

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-12-3	1721091-04	Water	08/01/17
DUP-4-3Q17	1721091-05	Water	08/01/17
MW-12-2	1721091-06	Water	08/01/17
MW-12-1	1721091-07	Water	08/01/17
MW-11-3	1721091-09	Water	08/01/17
MW-11-2	1721091-10	Water	08/01/17
DUP-5-3Q17	1721091-11	Water	08/01/17
MW-11-1	1721091-12	Water	08/01/17
EB-7-080117	1721091-13	Water	08/01/17
MW-12-3MS	1721091-04MS	Water	08/01/17
MW-12-3MSD	1721091-04MSD	Water	08/01/17
MW-12-3DUP	1721091-04DUP	Water	08/01/17

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following method:

Chromium by Environmental Protection Agency (EPA) Method 200.8

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. ICPMS Tune

The mass calibration was within 0.1 AMU and the percent relative standard deviation (%RSD) was less than or equal to 5%.

III. Instrument Calibration

Initial and continuing calibrations were performed as required by the method.

The initial calibration verification (ICV) and continuing calibration verification (CCV) standards were within QC limits.

IV. ICP Interference Check Sample Analysis

ICP interference check sample analyses were not required by the method.

V. Laboratory Blanks

Laboratory blanks were analyzed as required by the method. No contaminants were found in the laboratory blanks with the following exceptions:

Blank ID	Analyte	Maximum Concentration	Associated Samples
PB (prep blank)	Chromium	1.2170 ug/L	All samples in SDG 17-21091
ICB/CCB	Chromium	0.1700 ug/L	DUP-4-3Q17 MW-12-2 MW-12-1 MW-11-3 MW-11-2 DUP-5-3Q17 MW-11-1 EB-7-080117

Data qualification by the laboratory blanks was based on the maximum contaminant concentration in the laboratory blanks in the analysis of each analyte. The sample concentrations were either not detected or were significantly greater (>5X blank contaminants) than the concentrations found in the associated laboratory blanks with the following exceptions:

Sample	Analyte	Reported Concentration	Modified Final Concentration
MW-12-3	Chromium	1.8 ug/L	1.8U ug/L
DUP-4-3Q17	Chromium	2.4 ug/L	2.4U ug/L
MW-12-2	Chromium	1.9 ug/L	1.9U ug/L
MW-12-1	Chromium	2.9 ug/L	2.9U ug/L
MW-11-3	Chromium	2.9 ug/L	2.9U ug/L
MW-11-2	Chromium	2.1 ug/L	2.1U ug/L
DUP-5-3Q17	Chromium	2.0 ug/L	2.0U ug/L
MW-11-1	Chromium	2.2 ug/L	2.2U ug/L
EB-7-080117	Chromium	2.4 ug/L	2.4U ug/L

VI. Field Blanks

Sample EB-7-080117 was identified as an equipment blank. No contaminants were found with the following exceptions:

Blank ID	Analyte	Concentration
EB-7-080117	Chromium	2.4 ug/L

VII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VIII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

IX. Serial Dilution

Serial dilution was not performed for this SDG.

X. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the method. Percent recoveries (%R) were within QC limits.

XI. Field Duplicates

Samples MW-12-3 and DUP-4-3Q17 and samples MW-11-2 and DUP-5-3Q17 were identified as field duplicates. No results were detected in any of the samples with the following exceptions:

Analyte	Concentration (ug/L)		RPD
	MW-12-3	DUP-4-3Q17	
Chromium	1.8	2.4	29

Analyte	Concentration (ug/L)		RPD
	MW-11-2	DUP-5-3Q17	
Chromium	2.1	2.0	5

XII. Internal Standards (ICP-MS)

Raw data were not reviewed for Level III validation.

XIII. Sample Result Verification

Raw data were not reviewed for Level III validation.

XIV. Overall Assessment of Data

The analysis was conducted within all specifications of the method. No results were rejected in this SDG.

Due to laboratory blank contamination, data were qualified as not detected in nine samples.

The quality control criteria reviewed, other than those discussed above, were met and are considered acceptable. Based upon the data validation all other results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Chromium - Data Qualification Summary - SDG 17-21091

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017
Chromium - Laboratory Blank Data Qualification Summary - SDG 17-21091

Sample	Analyte	Modified Final Concentration	A or P
MW-12-3	Chromium	1.8U ug/L	A
DUP-4-3Q17	Chromium	2.4U ug/L	A
MW-12-2	Chromium	1.9U ug/L	A
MW-12-1	Chromium	2.9U ug/L	A
MW-11-3	Chromium	2.9U ug/L	A
MW-11-2	Chromium	2.1U ug/L	A
DUP-5-3Q17	Chromium	2.0U ug/L	A
MW-11-1	Chromium	2.2U ug/L	A
EB-7-080117	Chromium	2.4U ug/L	A



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:54:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-12-3

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-04

File ID: PE-EL3 170803R-181

Sampled: 08/01/17 08:15

Prepared: 08/03/17 11:00

Analyzed: 08/04/17 03:12

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0281

Sequence:

1713866

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.8 <i>U</i>	1	J	EPA-200.8

12/9/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:54:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

DUP-4-3Q17

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-05

File ID: PE-EL3 170803R-189

Sampled: 08/01/17 08:25

Prepared: 08/03/17 11:00

Analyzed: 08/04/17 03:39

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0281

Sequence:

1713866

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.4 U	1	J	EPA-200.8

Handwritten signature/initials



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:54:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-12-2

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-06

File ID: PE-EL3_170803R-190

Sampled: 08/01/17 09:15

Prepared: 08/03/17 11:00

Analyzed: 08/04/17 03:43

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0281

Sequence:

1713866

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	1.9 u	1	J	EPA-200.8

8092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:54:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

MW-12-1

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-07

File ID: PE-EL3 170803R-191

Sampled: 08/01/17 09:45

Prepared: 08/03/17 11:00

Analyzed: 08/04/17 03:46

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0281

Sequence:

1713866

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.9 U	1	J	EPA-200.8

209217



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:54:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-11-3

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-09

File ID: PE-EL3 170803R-192

Sampled: 08/01/17 11:30

Prepared: 08/03/17 11:00

Analyzed: 08/04/17 03:50

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0281

Sequence:

1713866

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.9 U	1	J	EPA-200.8

2092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:54:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-11-2

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-10

File ID: PE-EL3 170803R-193

Sampled: 08/01/17 11:50

Prepared: 08/03/17 11:00

Analyzed: 08/04/17 03:53

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0281

Sequence:

1713866

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.1 <i>✓</i>	1	J	EPA-200.8

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:54:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

DUP-5-3Q17

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-11

File ID: PE-EL3 170803R-194

Sampled: 08/01/17 12:00

Prepared: 08/03/17 11:00

Analyzed: 08/04/17 03:56

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BH0281

Sequence:

1713866

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.0 U	1	J	EPA-200.8

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:54:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-11-1

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-12

File ID: PE-EL3 170803R-195

Sampled: 08/01/17 12:50

Prepared: 08/03/17 11:00

Analyzed: 08/04/17 04:00

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0281

Sequence:

1713866

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.2 U	1	J	EPA-200.8

2092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:54:52AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

EB-7-080117

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-13

File ID: PE-EL3 170803R-196

Sampled: 08/01/17 13:00

Prepared: 08/03/17 11:00

Analyzed: 08/04/17 04:03

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0281

Sequence:

1713866

Calibration: UNASSIGNED

Instrument: PE-EL3

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	2.4 U	1	J	EPA-200.8

Handwritten signature/initials

LDC #: 39394A4a

VALIDATION COMPLETENESS WORKSHEET

Date: 09/15/17

SDG #: 17-21091

Level III

Page: 1 of 1

Laboratory: BC Laboratories, Inc.

Reviewer: ATV2nd Reviewer: B**METHOD:** Chromium (EPA Method 200.8)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A, A	
II.	ICP/MS Tune	A	
III.	Instrument Calibration	A	
IV.	ICP Interference Check Sample (ICS) Analysis	N	
V.	Laboratory Blanks	SW	
VI.	Field Blanks	SW	q = EB
VII.	Matrix Spike/Matrix Spike Duplicates	A	
VIII.	Duplicate sample analysis	A	
IX.	Serial Dilution	N	
X.	Laboratory control samples	A	ICS
XI.	Field Duplicates	SW	(1,2), (6,7)
XII.	Internal Standard (ICP-MS)	N	
XIII.	Sample Result Verification	N	
XIV.	Overall Assessment of Data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	MW-12-3	1721091-04	Water	08/01/17
2	DUP-4-3Q17	1721091-05	Water	08/01/17
3	MW-12-2	1721091-06	Water	08/01/17
4	MW-12-1	1721091-07	Water	08/01/17
5	MW-11-3	1721091-09	Water	08/01/17
6	MW-11-2	1721091-10	Water	08/01/17
7	DUP-5-3Q17	1721091-11	Water	08/01/17
8	MW-11-1	1721091-12	Water	08/01/17
9	EB-7-080117	1721091-13	Water	08/01/17
10	MW-12-3MS	1721091-04MS	Water	08/01/17
11	MW-12-3MSD	1721091-04MSD	Water	08/01/17
12	MW-12-3DUP	1721091-04DUP	Water	08/01/17
13				
14				

Notes:

VALIDATION FINDINGS WORKSHEET
PB/ICB/CCB QUALIFIED SAMPLES

METHOD: Trace metals (EPA SW 864 Method 6010B/6020/7000)
 Sample Concentration units, unless otherwise noted: ug/L

Soil preparation factor applied: NA
 Associated Samples: All

Analyte	Maximum PB ^a (mg/Kg)	Maximum PB ^a (ug/L)	Maximum ICB/CCB ^a (ug/L)	Action Level	1	2	3	4	5	6	7	8	9
Cr		1.2170		6.085	1.8	2.4	1.9	2.9	2.9	2.1	2.0	2.2	2.4

Sample Concentration units, unless otherwise noted: ug/L Associated Samples: 2 to 9

Analyte	Maximum PB ^a (mg/Kg)	Maximum PB ^a (ug/L)	Maximum ICB/CCB ^a (ug/L)	Action Level									
Cr			0.1700	0.85									

Samples with analyte concentrations within five times the associated ICB, CCB or PB concentration are listed above with the identifications from the Validation Completeness Worksheet. These sample results were qualified as not detected, "U".

Note : a - The listed analyte concentration is the highest ICB, CCB, or PB detected in the analysis of each element.

VALIDATION FINDINGS WORKSHEET
Field Blanks

METHOD: Trace Metals (EPA SW846 6010B/7000)

Blank units: ug/L **Associated sample units:** ug/L

Sampling date: 08/01/17 Soil factor applied _____

Field blank type: (circle one) Field Blank / Rinsate / Other: EB Associated Samples: X report

Analyte	Blank ID	Sample Identification												
		Action Limit												
Cr	2.4	12 U ³												

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:
 Samples with analyte concentrations within five times the associated field blank concentration are listed above, these sample results were qualified as not detected, "U".

LDC#: 39394A4a

VALIDATION FINDINGS WORKSHEET
Field Duplicates

Page: 1 of 1
Reviewer: ATL
2nd Reviewer: JL

METHOD: Metals (EPA Method 6010B/6020A7000)

Analyte	Concentration (ug/L)		RPD	
	1	2		
Chromium	1.8	2.4	29	

V:\FIELD DUPLICATES\Field Duplicates\FD_inorganic\2017\39394A4a.wpd

Analyte	Concentration (ug/L)		RPD	
	6	7		
Chromium	2.1	2.0	5	

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017
LDC Report Date: September 21, 2017
Parameters: Wet Chemistry
Validation Level: Level III & IV
Laboratory: BC Laboratories, Inc.
Sample Delivery Group (SDG): 17-21091

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-12-5	1721091-02	Water	08/01/17
MW-12-4**	1721091-03**	Water	08/01/17
MW-12-3	1721091-04	Water	08/01/17
DUP-4-3Q17	1721091-05	Water	08/01/17
MW-12-2	1721091-06	Water	08/01/17
MW-12-1	1721091-07	Water	08/01/17
MW-11-4	1721091-08	Water	08/01/17
MW-11-3	1721091-09	Water	08/01/17
MW-11-2	1721091-10	Water	08/01/17
DUP-5-3Q17	1721091-11	Water	08/01/17
MW-11-1	1721091-12	Water	08/01/17
EB-7-080117	1721091-13	Water	08/01/17
MW-12-5MS	1721091-02MS	Water	08/01/17
MW-12-5MSD	1721091-02MSD	Water	08/01/17
MW-12-5DUP	1721091-02DUP	Water	08/01/17
MW-12-3MS	1721091-04MS	Water	08/01/17
MW-12-3MSD	1721091-04MSD	Water	08/01/17
MW-12-3DUP	1721091-04DUP	Water	08/01/17
MW-11-1MS	1721091-12MS	Water	08/01/17
MW-11-1MSD	1721091-12MSD	Water	08/01/17
MW-11-1DUP	1721091-12DUP	Water	08/01/17

**Indicates sample underwent Level IV validation

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following methods:

Chloride, Nitrate as Nitrogen, and Sulfate by Environmental Protection Agency (EPA) Method 300.0

Nitrite as Nitrogen by EPA 353.2

Hexavalent Chromium by EPA SW 846 Method 7196

Orthophosphate as Phosphorus by EPA Method 365.1

Perchlorate by EPA Method 314.0

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results. Samples appended with a double asterisk on the cover page were subjected to Level IV data validation, which is comprised of the QC summary forms as well as the raw data, to confirm sample quantitation and identification.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. Initial Calibration

All criteria for the initial calibration of each method were met.

III. Continuing Calibration

Continuing calibration frequency and analysis criteria were met for each method when applicable.

IV. Laboratory Blanks

Laboratory blanks were analyzed as required by the methods. No contaminants were found in the laboratory blanks with the following exceptions:

Blank ID	Analyte	Maximum Concentration	Associated Samples
PB (prep blank)	Nitrite as Nitrogen	0.012671 mg/L	MW-11-1
ICB/CCB	Nitrite as Nitrogen	0.013528 mg/L	MW-11-1

Data qualification by the laboratory blanks was based on the maximum contaminant concentration in the laboratory blanks in the analysis of each analyte. The sample concentrations were either not detected or were significantly greater (>5X blank contaminants) than the concentrations found in the associated laboratory blanks with the following exceptions:

Sample	Analyte	Reported Concentration	Modified Final Concentration
MW-11-1	Nitrite as Nitrogen	0.022 mg/L	0.022U mg/L

V. Field Blanks

Sample EB-7-080117 was identified as an equipment blank. No contaminants were found.

VI. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

VIII. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the methods. Percent recoveries (%R) were within QC limits.

IX. Field Duplicates

Samples MW-12-3 and DUP-4-3Q17 and samples MW-11-2 and DUP-5-3Q17 were identified as field duplicates. No results were detected in any of the samples with the following exceptions:

Analyte	Concentration (ug/L)		RPD
	MW-12-3	DUP-4-2Q17	
Perchlorate	3.2	2.9	10

X. Sample Result Verification

All sample result verifications were acceptable for samples which underwent Level IV validation. Raw data were not reviewed for Level III validation.

XI. Overall Assessment of Data

The analysis was conducted within all specifications of the methods. No results were rejected in this SDG.

Due to laboratory blank contamination, data were qualified as not detected in one sample.

The quality control criteria reviewed, other than those discussed above, were met and are considered acceptable. Based upon the data validation all other results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Wet Chemistry - Data Qualification Summary - SDG 17-21091

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017
Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 17-21091

Sample	Compound	Modified Final Concentration	A or P
MW-11-1	Nitrite as Nitrogen	0.022U mg/L	A



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:11:15AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-300.0

MW-11-1

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-12

File ID: E080117.seq-45

Sampled: 08/01/17 12:50

Prepared: 08/01/17 18:00

Analyzed: 08/02/17 07:50

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH0197

Sequence:

1714072

Calibration: UNASSIGNED

Instrument: IC5

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
16887-00-6	Chloride	21	1		EPA-300.0
14797-55-8	Nitrate as N	0.53	1		EPA-300.0
14808-79-8	Sulfate	41	1		EPA-300.0

08/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:11:15AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-353.2

MW-11-1

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-12

File ID: 170802 0752 NO2-025

Sampled: 08/01/17 12:50

Prepared: 08/02/17 07:52

Analyzed: 08/02/17 08:14

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0292

Sequence:

1713792

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
14797-65-0	Nitrite as N	0.022 u	1	J	EPA-353.2

16092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:11:15AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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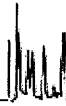
INORGANIC ANALYSIS DATA SHEET
EPA-365.1

MW-11-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-12</u>	File ID: <u>170802 0839 PO4-045</u>	
Sampled: <u>08/01/17 12:50</u>	Prepared: <u>08/02/17 08:39</u>	Analyzed: <u>08/02/17 08:39</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B/H0288</u>	Sequence: <u>1713795</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
---	ortho-Phosphate as P	0.017	1	U	EPA-365.1

top117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:11:15AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-12-3

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-04

File ID: 170801 2142 CR6-005

Sampled: 08/01/17 08:15

Prepared: 08/01/17 21:42

Analyzed: 08/01/17 21:42

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H0140

Sequence:

1713713

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

209217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:11:15AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-7196

DUP-4-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-05</u>	File ID: <u>170801 2142 CR6-009</u>	
Sampled: <u>08/01/17 08:25</u>	Prepared: <u>08/01/17 21:42</u>	Analyzed: <u>08/01/17 21:42</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B[H0140</u>	Sequence: <u>1713713</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>KONE-1</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

Handwritten signature: 8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:11:15AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-7196

MW-12-2

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-06

File ID: 170801 2142 CR6-028

Sampled: 08/01/17 09:15

Prepared: 08/01/17 21:42

Analyzed: 08/01/17 22:41

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BJH0140

Sequence:

1713713

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0014	2	UD	EPA-7196

SL892117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:11:15AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-12-1

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-07

File ID: 170801 2142 CR6-011

Sampled: 08/01/17 09:45

Prepared: 08/01/17 21:42

Analyzed: 08/01/17 21:43

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H0140

Sequence:

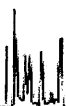
1713713

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:11:15AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-7196

MW-11-3

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-09

File ID: 170801 2142 CR6-012

Sampled: 08/01/17 11:30

Prepared: 08/01/17 21:42

Analyzed: 08/01/17 21:43

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B/H0140

Sequence: 1713713

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

2092117



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:11:15AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-7196

MW-11-2

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-10

File ID: 170801 2142 CR6-015

Sampled: 08/01/17 11:50

Prepared: 08/01/17 21:42

Analyzed: 08/01/17 21:46

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H0140

Sequence: 1713713

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:11:15AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

DUP-5-3Q17

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-11

File ID: 170801 2142 CR6-016

Sampled: 08/01/17 12:00

Prepared: 08/01/17 21:42

Analyzed: 08/01/17 21:46

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0140

Sequence:

1713713

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

209211



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:11:15AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-11-1

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-12

File ID: 170801 2142 CR6-017

Sampled: 08/01/17 12:50

Prepared: 08/01/17 21:42

Analyzed: 08/01/17 21:46

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H0140

Sequence:

1713713

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

809217



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:11:15AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-7196

EB-7-080117

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-13

File ID: 170801 2142 CR6-018

Sampled: 08/01/17 13:00

Prepared: 08/01/17 21:42

Analyzed: 08/01/17 21:46

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H0140

Sequence: 1713713

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8-09-2017



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:13:18AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-12-5

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-02

File ID: F081117.seq-11.0000.txt

Sampled: 08/01/17 07:30

Prepared: 08/11/17 14:00

Analyzed: 08/11/17 16:40

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH1247

Sequence:

1714506

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.7	1	J	EPA-314.0

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:13:18AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-12-4

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-03

File ID: F081117.seq-15.0000.txt

Sampled: 08/01/17 08:00

Prepared: 08/11/17 14:00

Analyzed: 08/11/17 17:36

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH1247

Sequence:

1714506

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.6	1	J	EPA-314.0

8-31-17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:13:18AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-314.0

MW-12-3

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-04

File ID: F081117.seq-16.0000.txt

Sampled: 08/01/17 08:15

Prepared: 08/11/17 14:00

Analyzed: 08/11/17 17:49

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH1247

Sequence:

1714506

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	3.2	1	J	EPA-314.0

8/29/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:13:18AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

DUP-4-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-05</u>	File ID: <u>F081117.seq-19.0000.txt</u>	
Sampled: <u>08/01/17 08:25</u>	Prepared: <u>08/11/17 14:00</u>	Analyzed: <u>08/11/17 18:31</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B H1247</u>	Sequence: <u>1714506</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.9	1	J	EPA-314.0

8092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:13:18AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-12-2

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-06

File ID: F081117.seq-20.0000.txt

Sampled: 08/01/17 09:15

Prepared: 08/11/17 14:00

Analyzed: 08/11/17 18:45

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H]1247

Sequence: 1714506

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:13:18AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-12-1

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-07

File ID: F081117.seq-21.0000.txt

Sampled: 08/01/17 09:45

Prepared: 08/11/17 14:00

Analyzed: 08/11/17 18:59

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H]1247

Sequence: 1714506

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:13:18AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET

EPA-314.0

MW-11-4

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-08

File ID: F081117.seq-22.0000.txt

Sampled: 08/01/17 11:00

Prepared: 08/11/17 14:00

Analyzed: 08/11/17 19:13

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH1247

Sequence: 1714506

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.94	1	J	EPA-314.0

8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:13:18AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-11-3

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-09</u>	File ID: <u>F081117.seq-23.0000.txt</u>	
Sampled: <u>08/01/17 11:30</u>	Prepared: <u>08/11/17 14:00</u>	Analyzed: <u>08/11/17 19:26</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B[H1247</u>	Sequence: <u>1714506</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

8/30/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:13:18AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-11-2

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-10

File ID: F081117.seq-24.0000.txt

Sampled: 08/01/17 11:50

Prepared: 08/11/17 14:00

Analyzed: 08/11/17 19:40

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH1247

Sequence:

1714506

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

SL08P17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:13:18AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

DUP-5-3Q17

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-11</u>	File ID: <u>F081117.seq-25.0000.txt</u>	
Sampled: <u>08/01/17 12:00</u>	Prepared: <u>08/11/17 14:00</u>	Analyzed: <u>08/11/17 19:54</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B[H]247</u>	Sequence: <u>1714506</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:13:18AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-11-1

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21091</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721091-12</u>	File ID: <u>F081117.seq-26.0000.txt</u>	
Sampled: <u>08/01/17 12:50</u>	Prepared: <u>08/11/17 14:00</u>	Analyzed: <u>08/11/17 20:08</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B H1247</u>	Sequence: <u>1714506</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.71	1	J	EPA-314.0

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:13:18AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

EB-7-080117

Laboratory: BC Laboratories

SDG: 17-21091

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721091-13

File ID: F081117.seq-27.0000.txt

Sampled: 08/01/17 13:00

Prepared: 08/11/17 14:00

Analyzed: 08/11/17 20:22

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B/H1247

Sequence: 1714506

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

Scop 117

LDC #: 39394A6

VALIDATION COMPLETENESS WORKSHEET

Date: 9/15/17

SDG #: 17-21091

Level III/IV

Page: 1 of 2

Laboratory: BC Laboratories, Inc.

Reviewer: ATL

2nd Reviewer: JB

METHOD: (Analyte) Chloride, Nitrate-N, Sulfate (EPA Method 300.0), Nitrite-N (EPA Method 353.2), Hexavalent Chromium (EPA SW846 Method 7196), Orthophosphate-P (EPA Method 365.1), Perchlorate (EPA Method 314.0).

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets:

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A, A	
II	Initial calibration	A	
III.	Calibration verification	A	
IV	Laboratory Blanks	SW	
V	Field blanks	ND	12 = EB
VI.	Matrix Spike/Matrix Spike Duplicates	A	
VII.	Duplicate sample analysis	A	
VIII.	Laboratory control samples	A	LCS
IX.	Field duplicates	SW	(3,4), (9,10)
X.	Sample result verification	A	Not reviewed for Level III validation.
XI	Overall assessment of data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

** Indicates sample was underwent Level IV review

	Client ID	Lab ID	Matrix	Date
1	MW-12-5	1721091-02	Water	08/01/17
2	MW-12-4**	1721091-03**	Water	08/01/17
3	MW-12-3	1721091-04	Water	08/01/17
4	DUP-4-3Q17	1721091-05	Water	08/01/17
5	MW-12-2	1721091-06	Water	08/01/17
6	MW-12-1	1721091-07	Water	08/01/17
7	MW-11-4	1721091-08	Water	08/01/17
8	MW-11-3	1721091-09	Water	08/01/17
9	MW-11-2	1721091-10	Water	08/01/17
10	DUP-5-3Q17	1721091-11	Water	08/01/17
11	MW-11-1	1721091-12	Water	08/01/17
12	EB-7-080117	1721091-13	Water	08/01/17
13	MW-12-5MS	1721091-02MS	Water	08/01/17
14	MW-12-5MSD	1721091-02MSD	Water	08/01/17
15	MW-12-5DUP	1721091-02DUP	Water	08/01/17
16	MW-12-3MS	1721091-04MS	Water	08/01/17
17	MW-12-3MSD	1721091-04MSD	Water	08/01/17

LDC #: 39394A6

VALIDATION COMPLETENESS WORKSHEET

Date: 09/15/17

SDG #: 17-21091

Level III/IV

Page: 2 of 2

Laboratory: BC Laboratories, Inc.

Reviewer: ATL

2nd Reviewer: JS

METHOD: (Analyte) Chloride, Nitrate-N, Sulfate (EPA Method 300.0), Nitrite-N (EPA Method 353.2), Hexavalent Chromium (EPA SW846 Method 7196), Orthophosphate-P (EPA Method 365.1), Perchlorate (EPA Method 314.0).

	Client ID	Lab ID	Matrix	Date
18	MW-12-3DUP	1721091-04DUP	Water	08/01/17
19	MW-11-1MS	1721091-12MS	Water	08/01/17
20	MW-11-1MSD	1721091-12MSD	Water	08/01/17
21	MW-11-1DUP	1721091-12DUP	Water	08/01/17
22				
23				
24				
25				
26				

Notes: _____

Method: Inorganics (EPA Method see cover)

Validation Area	Yes	No	NA	Findings/Comments
I. Technical holding times				
All technical holding times were met.	✓			
II. Calibration				
Were all instruments calibrated daily, each set-up time?		✓		
Were the proper number of standards used?	✓			
Were all initial calibration correlation coefficients > 0.995?	✓			
Were all initial and continuing calibration verification %Rs within the 90-110% QC limits?	✓			
Were titrant checks performed as required? (Level IV only)			✓	
Were balance checks performed as required? (Level IV only)			✓	
III. Blanks				
Was a method blank associated with every sample in this SDG?	✓			
Was there contamination in the method blanks? If yes, please see the Blanks validation completeness worksheet.	✓			
IV. Matrix spike/Matrix spike duplicates and Duplicates				
Were a matrix spike (MS) and duplicate (DUP) analyzed for each matrix in this SDG? If no, indicate which matrix does not have an associated MS/MSD or MS/DUP. Soil / Water.	✓			
Were the MS/MSD percent recoveries (%R) and the relative percent differences (RPD) within the 75-125 QC limits? If the sample concentration exceeded the spike concentration by a factor of 4 or more, no action was taken.	✓			
Were the MS/MSD or duplicate relative percent differences (RPD) ≤ 20% for waters and ≤ 35% for soil samples? A control limit of ≤ CRDL (≤ 2X CRDL for soil) was used for samples that were ≤ 5X the CRDL, including when only one of the duplicate sample values were < 5X the CRDL.	✓			
V. Laboratory control samples				
Was an LCS analyzed for this SDG?	✓			
Was an LCS analyzed per extraction batch?	✓			
Were the LCS percent recoveries (%R) and relative percent difference (RPD) within the 80-120% (85-115% for Method 300.0) QC limits?	✓			
VI. Regional Quality Assurance and Quality Control				
Were performance evaluation (PE) samples performed?		✓		
Were the performance evaluation (PE) samples within the acceptance limits?		✓		

VALIDATION FINDINGS CHECKLIST

Validation Area	Yes	No	NA	Findings/Comments
VII. Sample Result Verification				
Were RLs adjusted to reflect all sample dilutions and dry weight factors applicable to level IV validation?	✓			
Were detection limits < RL?	✓			
VIII. Overall assessment of data				
Overall assessment of data was found to be acceptable.	✓			
IX. Field duplicates				
Field duplicate pairs were identified in this SDG.	✓			
Target analytes were detected in the field duplicates.	✓			
X. Field blanks				
Field blanks were identified in this SDG.	✓			
Target analytes were detected in the field blanks.		✓		

VALIDATION FINDINGS WORKSHEET
Sample Specific Analysis Reference

All circled methods are applicable to each sample.

Sample ID	Parameter
11	pH TDS (Cl) F (N) (N) (SO ₄) (O-PO ₄) Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
3→6, 8→12	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC (Cr6+) ClO ₄
1→12	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ (ClO ₄)
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
QC	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
19, 20, 21	pH TDS Cl F NO ₃ (N) SO ₄ (O-PO ₄) Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
16, 17, 18	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC (Cr6+) ClO ₄
13, 14, 15	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ (ClO ₄)
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄

Comments: _____

VALIDATION FINDINGS WORKSHEET

Blanks

METHOD: Inorganics, Method See Cover

Conc. units: mg/L

Associated Samples: 11

Analyte	Blank ID	Blank ID	Blank Action Limit												
	PB	ICB/CCB (mg/L)		11											
Nitrite as N	0.012671		0.063355	0.022											
Nitrite as N		0.013528	0.06764	see above											

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:
 All contaminants within five times the method blank concentration were qualified as not detected, "U".

LDC# 39394A6

VALIDATION FINDINGS WORKSHEET
Field Duplicates

Page: 1 of 1
Reviewer: ATL
2nd Reviewer: SL

Inorganics: Method See Cover

Analyte	Concentration (ug/L)		RPD	
	3	4		
Perchlorate	3.2	2.9	10	

V:\FIELD DUPLICATES\Field Duplicates\FD_inorganic\2017\39394A6.wpd

LDC #: 39394AG

Validation Findings Worksheet
Initial and Continuing Calibration Calculation Verification

Page: 1 of 1
 Reviewer: ATV
 2nd Reviewer: VB

Method: Inorganics, Method See Cover

The correlation coefficient (r) for the calibration of ClO₄⁻ was recalculated. Calibration date: 08/01/17

An initial or continuing calibration verification percent recovery (%R) was recalculated for each type of analysis using the following formula:

$$\%R = \frac{\text{Found} \times 100}{\text{True}}$$

Where,

Found = concentration of each analyte measured in the analysis of the ICV or CCV solution

True = concentration of each analyte in the ICV or CCV source

Type of analysis	Analyte	Standard	Conc. (mg/L)	Area	Recalculated	Reported	Acceptable (Y/N)
					r or r ²	r or r ²	
Initial calibration	ClO ₄ ⁻	s1	0.0	0.0001	0.99942	0.99986	Y
		s2	2	0.0024			
		s3	4	0.0044			
		s4	6	0.007			
		s5	10	0.0119			
		s6	20	0.0243			
ICV (8/11 @ 15:17) Calibration verification	ClO ₄ ⁻	FOUND 9.1773	TRUE 10.000		91.8	92.2	Y
CCV ₁ (8/11 @ 18:03) Calibration verification	ClO ₄ ⁻	10.8205	10.000		108.2	110	Y
CCV ₂ (8/11 @ 22:33) Calibration verification	ClO ₄ ⁻	9.1773	10.000		91.8	95.6	Y

Comments: Refer to Calibration Verification findings worksheet for list of qualifications and associated samples when reported results do not agree within 10.0% of the recalculated results.

LDC #: 39394AG

VALIDATION FINDINGS WORKSHEET
Level IV Recalculation Worksheet

Page: 1 of 1
Reviewer: ATL
2nd Reviewer: VB

METHOD: Inorganics, Method see cover

Percent recoveries (%R) for a laboratory control sample and a matrix spike sample were recalculated using the following formula:

$\%R = \frac{\text{Found}}{\text{True}} \times 100$ Where, Found = concentration of each analyte measured in the analysis of the sample. For the matrix spike calculation, Found = SSR (spiked sample result) - SR (sample result).
True = concentration of each analyte in the source.

A sample and duplicate relative percent difference (RPD) was recalculated using the following formula:

$RPD = \frac{|S-D|}{(S+D)/2} \times 100$ Where, S = Original sample concentration
D = Duplicate sample concentration

Sample ID	Type of Analysis	Element	(mg/L) Found / S (units)	(mg/L) True / D (units)	Recalculated	Reported	Acceptable (Y/N)
					%R / RPD	%R / RPD	
LCS	Laboratory control sample (8/11 @ 16:26)	ClO ₄ ⁻	9.133 9.1773	10	91.3	89.5	Y
MS 13	Matrix spike sample (8/11 @ 17:08)	ClO ₄ ⁻	(SSR-SR) 9.754	10.101	96.6	96.8	Y
MS/MSD 13/14	Duplicate sample	ClO ₄ ⁻	13.285	12.461	6.41	4.99	Y

Comments: _____

VALIDATION FINDINGS WORKSHEET
Sample Calculation Verification

METHOD: Inorganics, Method see cover

Please see qualifications below for all questions answered "N". Not applicable questions are identified as "N/A".

- Y N N/A Have results been reported and calculated correctly?
- Y N N/A Are results within the calibrated range of the instruments?
- Y N N/A Are all detection limits below the CRQL?

Compound (analyte) results for ClO₄⁻ reported with a positive detect were recalculated and verified using the following equation:

Concentration = A x 830.28416209 Recalculation: 0.00261 x 830.28416209 = 2.167

#	Sample ID	Analyte	Reported Concentration (µg/L)	Calculated Concentration (µg/L)	Acceptable (Y/N)
	<u>2</u>	<u>ClO₄⁻</u>	<u>2.6</u>	<u>2.2</u>	<u>Y</u>

Note: _____

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017
LDC Report Date: September 21, 2017
Parameters: Volatiles
Validation Level: Level III
Laboratory: BC Laboratories, Inc.
Sample Delivery Group (SDG): 17-21528

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
TB-9-080317	1721528-01	Water	08/03/17
MW-10	1721528-02	Water	08/03/17
MW-16	1721528-03	Water	08/03/17
MW-7	1721528-04	Water	08/03/17
EB-9-080317	1721528-05	Water	08/03/17
MW-10MS	1721528-02MS	Water	08/03/17
MW-10MSD	1721528-02MSD	Water	08/03/17
MW-16MS	1721528-03MS	Water	08/03/17
MW-16MSD	1721528-03MSD	Water	08/03/17

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Superfund Organic Methods Data Review (June 2008). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following method:

Volatile Organic Compounds (VOCs) by Environmental Protection Agency (EPA) Method 524.2

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition and cooler temperatures upon receipt met validation criteria.

All technical holding time requirements were met.

II. GC/MS Instrument Performance Check

Instrument performance was checked at 12 hour intervals.

All ion abundance requirements were met.

III. Initial Calibration and Initial Calibration Verification

An initial calibration was performed as required by the method.

For compounds where average relative response factors (RRFs) were utilized, the percent relative standard deviations (%RSD) were less than or equal to 20.0%.

In the case where the laboratory used a calibration curve to evaluate the compounds, all coefficients of determination (r^2) were greater than or equal to 0.990.

The percent differences (%D) of the initial calibration verification (ICV) standard were less than or equal to 30.0% for all compounds.

IV. Continuing Calibration

Continuing calibration was performed at the required frequencies.

The percent differences (%D) were less than or equal to 30.0% for all compounds with the following exceptions:

Date	Compound	%D	Associated Samples	Flag	A or P
08/08/17	Bromomethane Methyl iodide Pentachloroethane	32.6 (≤ 30) 52.1 (≤ 30) 36.2 (≤ 30)	TB-9-080317 MW-10 MW-7 EB-9-080317	UJ (all non-detects) UJ (all non-detects) UJ (all non-detects)	P
08/08/17	Methyl iodide Pentachloroethane	54.1 (≤ 30) 41.0 (≤ 30)	MW-16	UJ (all non-detects) UJ (all non-detects)	P

V. Laboratory Blanks

Laboratory blanks were analyzed as required by the method. No contaminants were found in the laboratory blanks.

VI. Field Blanks

Sample TB-9-080317 was identified as a trip blank. No contaminants were found.

Sample EB-9-080317 was identified as an equipment blank. No contaminants were found.

VII. Surrogates

Surrogates were added to all samples as required by the method. All surrogate recoveries (%R) were within QC limits.

VIII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

IX. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the method. Percent recoveries (%R) were within QC limits.

X. Field Duplicates

No field duplicates were identified in this SDG.

XI. Internal Standards

All internal standard areas and retention times were within QC limits.

XII. Compound Quantitation

Raw data were not reviewed for Level III validation.

XIII. Target Compound Identifications

Raw data were not reviewed for Level III validation.

XIV. System Performance

Raw data were not reviewed for Level III validation.

XV. Overall Assessment of Data

The analysis was conducted within all specifications of the method. No results were rejected in this SDG.

Due to continuing calibration %D, data were qualified as estimated in five samples.

The quality control criteria reviewed, other than those discussed above, were met and are considered acceptable. Sample results that were found to be estimated (J) are usable for limited purposes only. Based upon the data validation all other results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Volatiles - Data Qualification Summary - SDG 17-21528

Sample	Compound	Flag	A or P	Reason
TB-9-080317 MW-10 MW-7 EB-9-080317	Bromomethane Methyl iodide Pentachloroethane	UJ (all non-detects) UJ (all non-detects) UJ (all non-detects)	P	Continuing calibration (%D)
MW-16	Methyl iodide Pentachloroethane	UJ (all non-detects) UJ (all non-detects)	P	Continuing calibration (%D)

NASA JPL, 3Q2017
Volatiles - Laboratory Blank Data Qualification Summary - SDG 17-21528

No Sample Data Qualified in this SDG



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:18:16AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

TB-9-080317

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21528</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721528-01</u>	File ID:	<u>07AUG46.D</u>		
Sampled:	<u>08/03/17 07:00</u>	Prepared:	<u>08/07/17 09:48</u>	Analyzed:	<u>08/08/17 05:13</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>BIH0652</u>	Sequence:	<u>1713910</u>	Calibration:	<u>1708011</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/09/2017



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:18:16AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

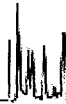
ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-10

Laboratory: BC Laboratories SDG: 17-21528
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721528-02 File ID: 07AUG39.D
Sampled: 08/03/17 07:15 Prepared: 08/07/17 09:48 Analyzed: 08/08/17 02:32
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B[H0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>UT</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

8/09/2017



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:18:16AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET

EPA-524.2

MW-10

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21528</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721528-02</u>	File ID:	<u>07AUG39.D</u>		
Sampled:	<u>08/03/17 07:15</u>	Prepared:	<u>08/07/17 09:48</u>	Analyzed:	<u>08/08/17 02:32</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>B[H0652</u>	Sequence:	<u>1713910</u>	Calibration:	<u>1708011</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:18:16AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

MW-16

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21528</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721528-03</u>	File ID:	<u>08AUG08.D</u>		
Sampled:	<u>08/03/17 08:30</u>	Prepared:	<u>08/08/17 07:00</u>	Analyzed:	<u>08/08/17 14:02</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>B[H0789</u>	Sequence:	<u>1714016</u>	Calibration:	<u>1708011</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

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Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:18:16AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET
EPA-524.2

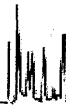
MW-16

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21528</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721528-03</u>	File ID:	<u>08AUG08.D</u>		
Sampled:	<u>08/03/17 08:30</u>	Prepared:	<u>08/08/17 07:00</u>	Analyzed:	<u>08/08/17 14:02</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>B[H0789</u>	Sequence:	<u>1714016</u>	Calibration:	<u>1708011</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

8292417



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:18:16AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-9-080317

Laboratory: BC Laboratories SDG: 17-21528
Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
Matrix: Water Laboratory ID: 1721528-05 File ID: 07AUG49.D
Sampled: 08/03/17 09:45 Prepared: 08/07/17 09:48 Analyzed: 08/08/17 06:22
Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
Batch: B/H0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
71-43-2	Benzene	1	0.11	U
108-86-1	Bromobenzene	1	0.15	U
74-97-5	Bromochloromethane	1	0.27	U
75-27-4	Bromodichloromethane	1	0.20	U
75-25-2	Bromoform	1	0.46	U
74-83-9	Bromomethane	1	0.20	U <i>UT</i>
104-51-8	n-Butylbenzene	1	0.15	U
135-98-8	sec-Butylbenzene	1	0.13	U
98-06-6	tert-Butylbenzene	1	0.18	U
56-23-5	Carbon tetrachloride	1	0.17	U
108-90-7	Chlorobenzene	1	0.14	U
75-00-3	Chloroethane	1	0.17	U
67-66-3	Chloroform	1	0.14	U
74-87-3	Chloromethane	1	0.11	U
95-49-8	2-Chlorotoluene	1	0.14	U
106-43-4	4-Chlorotoluene	1	0.093	U
124-48-1	Dibromochloromethane	1	0.22	U
96-12-8	1,2-Dibromo-3-chloropropane	1	0.89	U
106-93-4	1,2-Dibromoethane	1	0.22	U
74-95-3	Dibromomethane	1	0.23	U
95-50-1	1,2-Dichlorobenzene	1	0.21	U
541-73-1	1,3-Dichlorobenzene	1	0.16	U
106-46-7	1,4-Dichlorobenzene	1	0.15	U
75-71-8	Dichlorodifluoromethane	1	0.15	U
75-34-3	1,1-Dichloroethane	1	0.15	U
107-06-2	1,2-Dichloroethane	1	0.17	U
75-35-4	1,1-Dichloroethene	1	0.27	U
156-59-2	cis-1,2-Dichloroethene	1	0.27	U
156-60-5	trans-1,2-Dichloroethene	1	0.17	U
78-87-5	1,2-Dichloropropane	1	0.15	U
142-28-9	1,3-Dichloropropane	1	0.13	U
594-20-7	2,2-Dichloropropane	1	0.18	U

SL07-117



Tidewater Inc.
 3761 Attucks Drive
 Powell, OH 43065

Reported: 8/31/2017 11:18:16AM
 Project: JPL- GW Monitoring Wells
 Project Number: 3Q17
 Project Manager: David Conner

ORGANIC ANALYSIS DATA SHEET
EPA-524.2

EB-9-080317

Laboratory: BC Laboratories SDG: 17-21528
 Client: Tidewater Inc. Project: JPL- GW Monitoring Wells
 Matrix: Water Laboratory ID: 1721528-05 File ID: 07AUG49.D
 Sampled: 08/03/17 09:45 Prepared: 08/07/17 09:48 Analyzed: 08/08/17 06:22
 Solids: Preparation: EPA 5030 Water MS Initial/Final: 25 ml / 25 ml
 Batch: B[H0652 Sequence: 1713910 Calibration: 1708011 Instrument: MS-V5

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
563-58-6	1,1-Dichloropropene	1	0.19	U
10061-01-5	cis-1,3-Dichloropropene	1	0.14	U
10061-02-6	trans-1,3-Dichloropropene	1	0.13	U
100-41-4	Ethylbenzene	1	0.15	U
87-68-3	Hexachlorobutadiene	1	0.20	U
98-82-8	Isopropylbenzene	1	0.14	U
99-87-6	p-Isopropyltoluene	1	0.14	U
75-09-2	Methylene chloride	1	0.21	U
1634-04-4	Methyl t-butyl ether	1	0.14	U
91-20-3	Naphthalene	1	0.16	U
103-65-1	n-Propylbenzene	1	0.12	U
100-42-5	Styrene	1	0.12	U
630-20-6	1,1,1,2-Tetrachloroethane	1	0.21	U
79-34-5	1,1,2,2-Tetrachloroethane	1	0.17	U
127-18-4	Tetrachloroethene	1	0.23	U
108-88-3	Toluene	1	0.17	U
87-61-6	1,2,3-Trichlorobenzene	1	0.19	U
120-82-1	1,2,4-Trichlorobenzene	1	0.15	U
71-55-6	1,1,1-Trichloroethane	1	0.21	U
79-00-5	1,1,2-Trichloroethane	1	0.21	U
79-01-6	Trichloroethene	1	0.19	U
75-69-4	Trichlorofluoromethane	1	0.14	U
96-18-4	1,2,3-Trichloropropane	1	0.78	U
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	1	0.19	U
95-63-6	1,2,4-Trimethylbenzene	1	0.17	U
108-67-8	1,3,5-Trimethylbenzene	1	0.14	U
75-01-4	Vinyl chloride	1	0.18	U
67-64-1	Acetone	1	6.6	U
107-13-1	Acrylonitrile	1	1.5	U
107-05-1	Allyl chloride	1	0.47	U
994-05-8	t-Amyl Methyl ether	1	0.19	U
75-65-0	t-Butyl alcohol	1	9.4	U

8/31/17



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:18:16AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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ORGANIC ANALYSIS DATA SHEET

EPA-524.2

EB-9-080317

Laboratory:	<u>BC Laboratories</u>	SDG:	<u>17-21528</u>				
Client:	<u>Tidewater Inc.</u>	Project:	<u>JPL- GW Monitoring Wells</u>				
Matrix:	<u>Water</u>	Laboratory ID:	<u>1721528-05</u>	File ID:	<u>07AUG49.D</u>		
Sampled:	<u>08/03/17 09:45</u>	Prepared:	<u>08/07/17 09:48</u>	Analyzed:	<u>08/08/17 06:22</u>		
Solids:		Preparation:	<u>EPA 5030 Water MS</u>	Initial/Final:	<u>25 ml / 25 ml</u>		
Batch:	<u>B[H0652</u>	Sequence:	<u>1713910</u>	Calibration:	<u>1708011</u>	Instrument:	<u>MS-V5</u>

CAS NO.	COMPOUND	DILUTION	CONC. (ug/L)	Q
107-14-2	Chloroacetonitrile	1		U
109-69-3	1-Chlorobutane	1		U
513-88-2	1,1-Dichloropropanone	1		U
96-33-3	Methyl acrylate	1		U
98-95-3	Nitrobenzene	1		U
79-46-9	2-Nitropropane	1		U

* Values outside of QC limits

Signature

LDC #: 39394B1

VALIDATION COMPLETENESS WORKSHEET

SDG #: 17-21528

Level III

Laboratory: BC Laboratories, Inc.

Date: 9/12/17

Page: 1 of 1

Reviewer: [Signature]

2nd Reviewer: A

METHOD: GC/MS Volatiles (EPA Method 524.2)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A	
II.	GC/MS Instrument performance check	A	
III.	Initial calibration/ICV	A/A	RSD ≤ 20%. Y ² ICV ≤ 30%
IV.	Continuing calibration	M	CCV ≤ 30%
V.	Laboratory Blanks	A	
VI.	Field blanks	ND	TB=1. EB=5
VII.	Surrogate spikes	A	
VIII.	Matrix spike/Matrix spike duplicates	A	
IX.	Laboratory control samples	A	LCS
X.	Field duplicates	N	
XI.	Internal standards	A	
XII.	Compound quantitation RL/LOQ/LODs	N	
XIII.	Target compound identification	N	
XIV.	System performance	N	
XV.	Overall assessment of data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	TB-9-080317	1721528-01	Water	08/03/17
2	MW-10	1721528-02	Water	08/03/17
3	MW-16	1721528-03	Water	08/03/17
4	MW-7	1721528-04	Water	08/03/17
5	EB-9-080317	1721528-05	Water	08/03/17
6	MW-10MS	1721528-02MS	Water	08/03/17
7	MW-10MSD	1721528-02MSD	Water	08/03/17
8	MW-16MS	1721528-03MS	Water	08/03/17
9	MW-16MSD	1721528-03MSD	Water	08/03/17
10				
11				
12				
13				

TARGET COMPOUND WORKSHEET

METHOD: VOA

A. Chloromethane	AA. Tetrachloroethene	AAA. 1,3,5-Trimethylbenzene	AAAA. Ethyl tert-butyl ether	A1. 1,3-Butadiene
B. Bromomethane	BB. 1,1,2,2-Tetrachloroethane	BBB. 4-Chlorotoluene	BBBB. tert-Amyl methyl ether	B1. Hexane
C. Vinyl chloride	CC. Toluene	CCC. tert-Butylbenzene	CCCC. 1-Chlorohexane	C1. Heptane
D. Chloroethane	DD. Chlorobenzene	DDD. 1,2,4-Trimethylbenzene	DDDD. Isopropyl alcohol	D1. Propylene
E. Methylene chloride	EE. Ethylbenzene	EEE. sec-Butylbenzene	EEEE. Acetonitrile	E1. Freon 11
F. Acetone	FF. Styrene	FFF. 1,3-Dichlorobenzene	FFFF. Acrolein	F1. Freon 12
G. Carbon disulfide	GG. Xylenes, total	GGG. p-Isopropyltoluene	GGGG. Acrylonitrile	G1. Freon 113
H. 1,1-Dichloroethene	HH. Vinyl acetate	HHH. 1,4-Dichlorobenzene	HHHH. 1,4-Dioxane	H1. Freon 114
I. 1,1-Dichloroethane	II. 2-Chloroethylvinyl ether	III. n-Butylbenzene	IIII. Isobutyl alcohol	I1. 2-Nitropropane
J. 1,2-Dichloroethene, total	JJ. Dichlorodifluoromethane	JJJ. 1,2-Dichlorobenzene	JJJJ. Methacrylonitrile	J1. Dimethyl disulfide
K. Chloroform	KK. Trichlorofluoromethane	KKK. 1,2,4-Trichlorobenzene	KKKK. Propionitrile	K1. 2,3-Dimethyl pentane
L. 1,2-Dichloroethane	LL. Methyl-tert-butyl ether	LLL. Hexachlorobutadiene	LLLL. Ethyl ether	L1. 2,4-Dimethyl pentane
M. 2-Butanone	MM. 1,2-Dibromo-3-chloropropane	MMM. Naphthalene	MMMM. Benzyl chloride	M1. 3,3-Dimethyl pentane
N. 1,1,1-Trichloroethane	NN. Methyl ethyl ketone	NNN. 1,2,3-Trichlorobenzene	NNNN. Iodomethane	N1. 2-Methylpentane
O. Carbon tetrachloride	OO. 2,2-Dichloropropane	OOO. 1,3,5-Trichlorobenzene	OOOO. 1,1-Difluoroethane	O1. 3-Methylpentane
P. Bromodichloromethane	PP. Bromochloromethane	PPP. trans-1,2-Dichloroethene	PPPP. Tetrahydrofuran	P1. 3-Ethylpentane
Q. 1,2-Dichloropropane	QQ. 1,1-Dichloropropene	QQQ. cis-1,2-Dichloroethene	QQQQ. Methyl acetate	Q1. 2,2-Dimethylpentane
R. cis-1,3-Dichloropropene	RR. Dibromomethane	RRR. m,p-Xylenes	RRRR. Ethyl acetate	R1. 2,2,3-Trimethylbutane
S. Trichloroethene	SS. 1,3-Dichloropropane	SSS. o-Xylene	SSSS. Cyclohexane	S1. 2,2,4-Trimethylpentane
T. Dibromochloromethane	TT. 1,2-Dibromoethane	TTT. 1,1,2-Trichloro-1,2,2-trifluoroethane	TTTT. Methylcyclohexane	T1. 2-Methylhexane
U. 1,1,2-Trichloroethane	UU. 1,1,1,2-Tetrachloroethane	UUU. 1,2-Dichlorotetrafluoroethane	UUUU. Allyl chloride	U1. Nonanal
V. Benzene	VV. Isopropylbenzene	VVV. 4-Ethyltoluene	VVVV. Methyl methacrylate	V1. 2-Methylnaphthalene
W. trans-1,3-Dichloropropene	WW. Bromobenzene	WWW. Ethanol	WWWW. Ethyl methacrylate	W1. Methanol
X. Bromoform	XX. 1,2,3-Trichloropropane	XXX. Di-isopropyl ether	XXXX. cis-1,4-Dichloro-2-butene	X1. 1,2,3-Trimethylbenzene
Y. 4-Methyl-2-pentanone	YY. n-Propylbenzene	YYY. tert-Butanol	YYYY. trans-1,4-Dichloro-2-butene	Y1.
Z. 2-Hexanone	ZZ. 2-Chlorotoluene	ZZZ. tert-Butyl alcohol	ZZZZ. Pentachloroethane	Z1.

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017

LDC Report Date: September 20, 2017

Parameters: Chromium

Validation Level: Level III

Laboratory: BC Laboratories, Inc.

Sample Delivery Group (SDG): 17-21528

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-10	1721528-02	Water	08/03/17
MW-16	1721528-03	Water	08/03/17
MW-7	1721528-04	Water	08/03/17
EB-9-080317	1721528-05	Water	08/03/17
MW-10MS	1721528-02MS	Water	08/03/17
MW-10MSD	1721528-02MSD	Water	08/03/17
MW-10DUP	1721528-02DUP	Water	08/03/17

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following method:

Chromium by Environmental Protection Agency (EPA) Method 200.8

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. ICPMS Tune

The mass calibration was within 0.1 AMU and the percent relative standard deviation (%RSD) was less than or equal to 5%.

III. Instrument Calibration

Initial and continuing calibrations were performed as required by the method.

The initial calibration verification (ICV) and continuing calibration verification (CCV) standards were within QC limits.

IV. ICP Interference Check Sample Analysis

ICP interference check sample analyses were not required by the method.

V. Laboratory Blanks

Laboratory blanks were analyzed as required by the method. No contaminants were found in the laboratory blanks with the following exceptions:

Blank ID	Analyte	Maximum Concentration	Associated Samples
ICB/CCB	Chromium	1.6670 ug/L	MW-7

Data qualification by the laboratory blanks was based on the maximum contaminant concentration in the laboratory blanks in the analysis of each analyte. The sample concentrations were either not detected or were significantly greater (>5X blank contaminants) than the concentrations found in the associated laboratory blanks.

VI. Field Blanks

Sample EB-9-080317 was identified as an equipment blank. No contaminants were found with the following exceptions:

Blank ID	Analyte	Concentration
EB-9-080317	Chromium	0.78 ug/L

VII. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VIII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

IX. Serial Dilution

Serial dilution was not performed for this SDG.

X. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the method. Percent recoveries (%R) were within QC limits.

XI. Field Duplicates

No field duplicates were identified in this SDG.

XII. Internal Standards (ICP-MS)

Raw data were not reviewed for Level III validation.

XIII. Sample Result Verification

Raw data were not reviewed for Level III validation.

XIV. Overall Assessment of Data

The analysis was conducted within all specifications of the method. No results were rejected in this SDG.

The quality control criteria reviewed, were met and are considered acceptable. Based upon the data validation all results are considered valid and usable for all purposes.

NASA JPL, 3Q2017
Chromium - Data Qualification Summary - SDG 17-21528

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017
Chromium - Laboratory Blank Data Qualification Summary - SDG 17-21528

No Sample Data Qualified in this SDG



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:22:46AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-10

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-02

File ID: PE_EL2_170807-248

Sampled: 08/03/17 07:15

Prepared: 08/07/17 09:30

Analyzed: 08/07/17 23:59

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0588

Sequence:

1713993

Calibration: UNASSIGNED

Instrument: PE-EL2

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	3.8	1		EPA-200.8

Handwritten signature/initials



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:22:46AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-16

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-03

File ID: PE_EL2_170807-259

Sampled: 08/03/17 08:30

Prepared: 08/07/17 09:30

Analyzed: 08/08/17 00:38

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BIH0588

Sequence: 1713993

Calibration: UNASSIGNED

Instrument: PE-EL2

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	76	1		EPA-200.8

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:22:46AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-200.8

MW-7

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-04

File ID: PE_EL2_170808-225

Sampled: 08/03/17 09:35

Prepared: 08/07/17 09:30

Analyzed: 08/09/17 05:40

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: BJH0588

Sequence: 1714090

Calibration: UNASSIGNED

Instrument: PE-EL2

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	7400	10	D	EPA-200.8

7/29/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:22:46AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-200.8

EB-9-080317

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-05

File ID: PE_EL2_170807-261

Sampled: 08/03/17 09:45

Prepared: 08/07/17 09:30

Analyzed: 08/08/17 00:44

Solids: 0.00

Preparation: EPA 200.2

Initial/Final: 50 ml / 50 ml

Batch: B[H0588

Sequence: 1713993

Calibration: UNASSIGNED

Instrument: PE-EL2

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
7440-47-3	Total Recoverable Chromium	0.78	1	J	EPA-200.8

Handwritten signature

LDC #: 39394B4a

VALIDATION COMPLETENESS WORKSHEET

Date: 09/18/17

SDG #: 17-21528

Level III

Page: 1 of 1

Laboratory: BC Laboratories, Inc.

Reviewer: ATL
2nd Reviewer: JB

METHOD: Chromium (EPA Method 200.8)

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A, A	
II.	ICP/MS Tune	A	
III.	Instrument Calibration	A	
IV.	ICP Interference Check Sample (ICS) Analysis	N	
V.	Laboratory Blanks	SW	
VI.	Field Blanks	SW	4 = EB
VII.	Matrix Spike/Matrix Spike Duplicates	A	
VIII.	Duplicate sample analysis	A	
IX.	Serial Dilution	N	
X.	Laboratory control samples	A	LCS
XI.	Field Duplicates	N	
XII.	Internal Standard (ICP-MS)	N	
XIII.	Sample Result Verification	N	
XIV.	Overall Assessment of Data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	MW-10	1721528-02	Water	08/03/17
2	MW-16	1721528-03	Water	08/03/17
3	MW-7	1721528-04	Water	08/03/17
4	EB-9-080317	1721528-05	Water	08/03/17
5	MW-10MS	1721528-02MS	Water	08/03/17
6	MW-10MSD	1721528-02MSD	Water	08/03/17
7	MW-10DUP	1721528-02DUP	Water	08/03/17
8				
9				
10				
11				
12				
13				
14				

Notes:

VALIDATION FINDINGS WORKSHEET
PB/ICB/CCB QUALIFIED SAMPLES

METHOD: Trace metals (EPA SW 864 Method 6010B/6020/7000)

Soil preparation factor applied: NA

Sample Concentration units, unless otherwise noted: ug/L

Associated Samples: 3

Analyte	Maximum PB ^a (mg/Kg)	Maximum PB ^a (ug/L)	Maximum ICB/CCB ^a (ug/L)	Action Level									
Cr			1.6670	8.335									

Samples with analyte concentrations within five times the associated ICB, CCB or PB concentration are listed above with the identifications from the Validation Completeness Worksheet. These sample results were qualified as not detected, "U".

Note : a - The listed analyte concentration is the highest ICB, CCB, or PB detected in the analysis of each element.

VALIDATION FINDINGS WORKSHEET
Field Blanks

METHOD: Trace Metals (EPA SW846 6010B/7000)

Blank units: ug/L **Associated sample units:** ug/L

Sampling date: 08/03/17 Soil factor applied _____

Field blank type: (circle one) Field Blank / Rinsate / Other: EB Associated Samples: X report

Analyte	Blank ID	Sample Identification									
		Action Limit									
Cr	0.78	3.90									

CIRCLED RESULTS WERE NOT QUALIFIED. ALL RESULTS NOT CIRCLED WERE QUALIFIED BY THE FOLLOWING STATEMENT:
 Samples with analyte concentrations within five times the associated field blank concentration are listed above, these sample results were qualified as not detected, "U".

Laboratory Data Consultants, Inc. Data Validation Report

Project/Site Name: NASA JPL, 3Q2017
LDC Report Date: September 20, 2017
Parameters: Wet Chemistry
Validation Level: Level III
Laboratory: BC Laboratories, Inc.
Sample Delivery Group (SDG): 17-21528

Sample Identification	Laboratory Sample Identification	Matrix	Collection Date
MW-10	1721528-02	Water	08/03/17
MW-16	1721528-03	Water	08/03/17
MW-7	1721528-04	Water	08/03/17
EB-9-080317	1721528-05	Water	08/03/17
MW-10MS	1721528-02MS	Water	08/03/17
MW-10MSD	1721528-02MSD	Water	08/03/17
MW-10DUP	1721528-02DUP	Water	08/03/17
MW-16MS	1721528-03MS	Water	08/03/17
MW-16MSD	1721528-03MSD	Water	08/03/17
MW-16DUP	1721528-03DUP	Water	08/03/17

Introduction

This Data Validation Report (DVR) presents data validation findings and results for the associated samples listed on the cover page. Data validation was performed in accordance with a modified outline of the USEPA Contract Laboratory Program National Functional Guidelines (CLPNFG) for Inorganic Superfund Data Review (January 2010). Where specific guidance was not available, the data has been evaluated in a conservative manner consistent with industry standards using professional experience.

The analyses were performed by the following methods:

Chloride, Nitrate as Nitrogen, Sulfate Environmental Protection Agency (EPA) Method 300.0

Nitrite as Nitrogen EPA Method 353.2

Hexavalent Chromium by EPA SW 846 Method 7196

Orthophosphate-Phosphorus by EPA Method 365.1

Perchlorate by EPA Method 314.0

All sample results were subjected to Level III data validation, which comprises an evaluation of quality control (QC) summary results.

The following are definitions of the data qualifiers utilized during data validation:

- J (Estimated): The compound or analyte was analyzed for and positively identified by the laboratory; however the reported concentration is estimated due to non-conformances discovered during data validation.
- U (Non-detected): The compound or analyte was analyzed for and positively identified by the laboratory; however the compound or analyte should be considered non-detected at the reported concentration due to the presence of contaminants detected in the associated blank(s).
- UJ (Non-detected estimated): The compound or analyte was reported as not detected by the laboratory; however the reported quantitation/detection limit is estimated due to non-conformances discovered during data validation.
- R (Rejected): The sample results were rejected due to gross non-conformances discovered during data validation. Data qualified as rejected is not usable.
- NA (Not Applicable): The non-conformance discovered during data validation demonstrates a high bias, while the affected compound or analyte in the associated sample(s) was reported as not detected by the laboratory and did not warrant the qualification of the data.

A qualification summary table is provided at the end of this report if data has been qualified. Flags are classified as P (protocol) or A (advisory) to indicate whether the flag is due to a laboratory deviation from a specified protocol or is of technical advisory nature.

I. Sample Receipt and Technical Holding Times

All samples were received in good condition.

All technical holding time requirements were met.

II. Initial Calibration

All criteria for the initial calibration of each method were met.

III. Continuing Calibration

Continuing calibration frequency and analysis criteria were met for each method when applicable.

IV. Laboratory Blanks

Laboratory blanks were analyzed as required by the methods. No contaminants were found in the laboratory blanks.

V. Field Blanks

Sample EB-9-080317 was identified as an equipment blank. No contaminants were found.

VI. Matrix Spike/Matrix Spike Duplicates

Matrix spike (MS) and matrix spike duplicate (MSD) sample analysis was performed on an associated project sample. Percent recoveries (%R) were within QC limits. Relative percent differences (RPD) were within QC limits.

VII. Duplicate Sample Analysis

Duplicate (DUP) sample analysis was performed on an associated project sample. Results were within QC limits.

VIII. Laboratory Control Samples

Laboratory control samples (LCS) were analyzed as required by the methods. Percent recoveries (%R) were within QC limits.

IX. Field Duplicates

No field duplicates were identified in this SDG.

X. Sample Result Verification

Raw data were not reviewed for Level III validation.

XI. Overall Assessment of Data

The analysis was conducted within all specifications of the methods. No results were rejected in this SDG.

The quality control criteria reviewed were met and are considered acceptable. Based upon the data validation all results are considered valid and usable for all purposes.

NASA JPL, 3Q2017

Wet Chemistry - Data Qualification Summary - SDG 17-21528

No Sample Data Qualified in this SDG

NASA JPL, 3Q2017

Wet Chemistry - Laboratory Blank Data Qualification Summary - SDG 17-21528

No Sample Data Qualified in this SDG



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:20:22AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-300.0

MW-16

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21528</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721528-03</u>	File ID: <u>E080417.seq-20</u>	
Sampled: <u>08/03/17 08:30</u>	Prepared: <u>08/04/17 17:00</u>	Analyzed: <u>08/04/17 18:29</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B[H0502</u>	Sequence: <u>1714411</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC5</u>

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
16887-00-6	Chloride	86	1		EPA-300.0
14797-55-8	Nitrate as N	1.4	1		EPA-300.0
14808-79-8	Sulfate	54	1		EPA-300.0

2092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-300.0

MW-7

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-04

File ID: E080417.seq-21

Sampled: 08/03/17 09:35

Prepared: 08/04/17 17:00

Analyzed: 08/04/17 18:47

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H0502

Sequence: 1714411

Calibration: UNASSIGNED

Instrument: IC5

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
16887-00-6	Chloride	86	1		EPA-300.0
14797-55-8	Nitrate as N	1.5	1		EPA-300.0
14808-79-8	Sulfate	55	1		EPA-300.0

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-10

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-02

File ID: F081417.seq-49.0000.txt

Sampled: 08/03/17 07:15

Prepared: 08/15/17 01:00

Analyzed: 08/15/17 08:45

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B[H1387

Sequence: 1714797

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	2.0	1	J	EPA-314.0

10892117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-16

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-03

File ID: F081417.seq-97.0000.txt

Sampled: 08/03/17 08:30

Prepared: 08/15/17 01:00

Analyzed: 08/15/17 21:54

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: B/H1591

Sequence: 1714797

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-314.0

MW-7

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-04RE2

File ID: F081417.seq-105.0000.txt

Sampled: 08/03/17 09:35

Prepared: 08/15/17 01:00

Analyzed: 08/15/17 23:49

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH1387

Sequence: 1714797

Calibration: UNASSIGNED

Instrument: IC6

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	12	1		EPA-314.0

11/29/2017



Tidewater Inc. 3761 Attucks Drive Powell, OH 43065	Reported: 8/31/2017 11:20:22AM Project: JPL- GW Monitoring Wells Project Number: 3Q17 Project Manager: David Conner
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INORGANIC ANALYSIS DATA SHEET
EPA-314.0

EB-9-080317

Laboratory: <u>BC Laboratories</u>	SDG: <u>17-21528</u>		
Client: <u>Tidewater Inc.</u>	Project: <u>JPL- GW Monitoring Wells</u>		
Matrix: <u>Water</u>	Laboratory ID: <u>1721528-05</u>	File ID: <u>F081417.seq-62.0000.txt</u>	
Sampled: <u>08/03/17 09:45</u>	Prepared: <u>08/15/17 01:00</u>	Analyzed: <u>08/15/17 11:45</u>	
Solids: <u>0.00</u>	Preparation: <u>No Prep</u>	Initial/Final: <u>20 ml / 20 ml</u>	
Batch: <u>B/H1387</u>	Sequence: <u>1714797</u>	Calibration: <u>UNASSIGNED</u>	Instrument: <u>IC6</u>

CAS NO.	Analyte	Concentration (ug/L)	Dilution Factor	Q	Method
14797-73-0	Perchlorate	0.58	1	U	EPA-314.0

8/31/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-353.2

MW-16

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-03

File ID: 170804 0816 NO2-025

Sampled: 08/03/17 08:30

Prepared: 08/04/17 08:16

Analyzed: 08/04/17 08:39

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH0564

Sequence: 1713981

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
14797-65-0	Nitrite as N	0.010	1	U	EPA-353.2

2092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-353.2

MW-7

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-04

File ID: 170804 0816 NO2-029

Sampled: 08/03/17 09:35

Prepared: 08/04/17 08:16

Analyzed: 08/04/17 08:39

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0564

Sequence: 1713981

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
14797-65-0	Nitrite as N	0.010	1	U	EPA-353.2

10/2/17



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-365.1

MW-16

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-03

File ID: 170804 0909 PO4-045

Sampled: 08/03/17 08:30

Prepared: 08/04/17 09:09

Analyzed: 08/04/17 09:09

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BH0566

Sequence: 1713984

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
---	ortho-Phosphate as P	0.30	1		EPA-365.1

8/31/17



Tidewater Inc.
3761 Attucks Drive
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Reported: 8/31/2017 11:20:22AM
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Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-365.1

MW-7

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-04

File ID: 170804 0909 PO4-049

Sampled: 08/03/17 09:35

Prepared: 08/04/17 09:09

Analyzed: 08/04/17 09:09

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0566

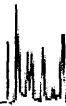
Sequence: 1713984

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
---	ortho-Phosphate as P	0.017	1	U	EPA-365.1

Handwritten signature/initials



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET

EPA-7196

MW-10

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-02

File ID: 170803 2110 CR6-005

Sampled: 08/03/17 07:15

Prepared: 08/03/17 21:10

Analyzed: 08/03/17 21:10

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0395

Sequence: 1713941

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0029	1		EPA-7196

2017



Tidewater Inc.
3761 Attucks Drive
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Reported: 8/31/2017 11:20:22AM
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Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-16

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-03

File ID: 170803 2110 CR6-011

Sampled: 08/03/17 08:30

Prepared: 08/03/17 21:10

Analyzed: 08/03/17 21:10

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0396

Sequence: 1713941

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0016	1	J	EPA-7196

8092117



Tidewater Inc.
3761 Attucks Drive
Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

MW-7

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-04

File ID: 170803 2110 CR6-023

Sampled: 08/03/17 09:35

Prepared: 08/03/17 21:10

Analyzed: 08/03/17 21:31

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0396

Sequence: 1713941

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.0011	1	J	EPA-7196

Handwritten signature/initials



Tidewater Inc.
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Powell, OH 43065

Reported: 8/31/2017 11:20:22AM
Project: JPL- GW Monitoring Wells
Project Number: 3Q17
Project Manager: David Conner

INORGANIC ANALYSIS DATA SHEET
EPA-7196

EB-9-080317

Laboratory: BC Laboratories

SDG: 17-21528

Client: Tidewater Inc.

Project: JPL- GW Monitoring Wells

Matrix: Water

Laboratory ID: 1721528-05

File ID: 170803 2110 CR6-018

Sampled: 08/03/17 09:45

Prepared: 08/03/17 21:10

Analyzed: 08/03/17 21:14

Solids: 0.00

Preparation: No Prep

Initial/Final: 20 ml / 20 ml

Batch: BIH0396

Sequence: 1713941

Calibration: UNASSIGNED

Instrument: KONE-1

CAS NO.	Analyte	Concentration (mg/L)	Dilution Factor	Q	Method
18540-29-9	Hexavalent Chromium	0.00070	1	U	EPA-7196

8092117

LDC #: 39394B6

VALIDATION COMPLETENESS WORKSHEET

Date: 09/18/17

SDG #: 17-21528

Level III

Page: 1 of 1

Laboratory: BC Laboratories, Inc.

Reviewer: AJL

2nd Reviewer: JB

METHOD: (Analyte) Chloride, Nitrate-N, Sulfate (EPA Method 300.0), Nitrite-N (EPA Method 353.2), Hexavalent Chromium (EPA SW846 Method 7196), Orthophosphate-P (EPA Method 365.1), Perchlorate (EPA Method 314.0).

The samples listed below were reviewed for each of the following validation areas. Validation findings are noted in attached validation findings worksheets.

	Validation Area		Comments
I.	Sample receipt/Technical holding times	A/A	
II	Initial calibration	A	
III.	Calibration verification	A	
IV	Laboratory Blanks	A	
V	Field blanks	ND	4 = EB
VI.	Matrix Spike/Matrix Spike Duplicates	A	
VII.	Duplicate sample analysis	A	
VIII.	Laboratory control samples	A	LCS
IX.	Field duplicates	N	
X.	Sample result verification	N	
XI	Overall assessment of data	A	

Note: A = Acceptable
N = Not provided/applicable
SW = See worksheet

ND = No compounds detected
R = Rinsate
FB = Field blank

D = Duplicate
TB = Trip blank
EB = Equipment blank

SB=Source blank
OTHER:

	Client ID	Lab ID	Matrix	Date
1	MW-10	1721528-02	Water	08/03/17
2	MW-16	1721528-03	Water	08/03/17
3	MW-7	1721528-04	Water	08/03/17
4	EB-9-080317	1721528-05	Water	08/03/17
5	MW-10MS	1721528-02MS	Water	08/03/17
6	MW-10MSD	1721528-02MSD	Water	08/03/17
7	MW-10DUP	1721528-02DUP	Water	08/03/17
8	MW-16MS	1721528-03MS	Water	08/03/17
9	MW-16MSD	1721528-03MSD	Water	08/03/17
10	MW-16DUP	1721528-03DUP	Water	08/03/17
11				
12				
13				
14				
15				

Notes:

VALIDATION FINDINGS WORKSHEET
Sample Specific Analysis Reference

All circled methods are applicable to each sample.

Sample ID	Parameter
2,3	pH TDS (Cl) F (N) (N) (SO ₄) (O-PO ₄) Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
1-4	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC (Cr6+) (ClO ₄)
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
QC	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
8,9,10	pH TDS (Cl) F (N) (N) (SO ₄) (O-PO ₄) Alk CN NH ₃ TKN TOC (Cr6+) (ClO ₄)
5,6,7	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC (Cr6+) (ClO ₄)
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄
	pH TDS Cl F NO ₃ NO ₂ SO ₄ O-PO ₄ Alk CN NH ₃ TKN TOC Cr6+ ClO ₄

Comments: _____