

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX111618\
 Data File : VX006055.D
 Acq On : 16 Nov 2018 17:35
 Operator : JC/MD
 Sample : J5965-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-7

Quant Time: Nov 17 02:44:24 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 07 14:40:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	136170	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	191189	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	182263	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	79335	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.07	65	80032	51.12	ug/l	0.00
Spiked Amount			Recovery	=	102.24%	
35) Dibromofluoromethane	5.51	113	65179	50.12	ug/l	0.00
Spiked Amount			Recovery	=	100.24%	
50) Toluene-d8	8.71	98	217706	50.36	ug/l	0.00
Spiked Amount			Recovery	=	100.72%	
62) 4-Bromofluorobenzene	11.14	95	73533	47.31	ug/l	0.00
Spiked Amount			Recovery	=	94.62%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Chloromethane	1.33	50	1469	0.912	ug/l	96
11) Tert butyl alcohol	3.06	59	31402	70.024	ug/l	99
16) Acetone	2.45	43	5047	5.988	ug/l	98
18) Methyl Acetate	2.78	43	7790	2.026	ug/l	100
22) Diisopropyl ether	3.87	45	1318	0.280	ug/l #	66
31) Cyclohexane	5.58	56	4704	2.157	ug/l #	95
39) Methylcyclohexane	7.46	83	1983	1.062	ug/l	85
40) Benzene	6.15	78	19040	3.544	ug/l	99
67) Ethyl Benzene	10.26	91	2662	0.455	ug/l #	87
73) Isopropylbenzene	11.02	105	2518	0.549	ug/l	99
78) n-propylbenzene	11.36	91	2222	0.411	ug/l	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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