

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX120318\
 Data File : VX006245.D
 Acq On : 03 Dec 2018 11:37
 Operator : JC/MD
 Sample : J6147-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-11

Quant Time: Dec 04 05:42:32 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X112918W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 30 03:55:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	963116	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	1412186	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	1274321	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	624461	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.07	65	463363	47.13	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.26%	
35) Dibromofluoromethane	5.51	113	452868	49.53	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.06%	
50) Toluene-d8	8.71	98	1642253	49.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.76%	
62) 4-Bromofluorobenzene	11.13	95	603389	48.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.30%	
Target Compounds						
16) Acetone	2.45	43	22676	4.915	ug/l	98
17) Carbon Disulfide	2.57	76	17934	0.668	ug/l	98
27) cis-1,2-Dichloroethene	4.61	96	16921	1.633	ug/l	92
30) Chloroform	5.21	83	6477	0.401	ug/l	94
44) Trichloroethene	7.21	130	15578	1.444	ug/l	86
64) Tetrachloroethene	9.34	164	179516	18.706	ug/l	96

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

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