

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX032720\
 Data File : VX015419.D
 Acq On : 28 Mar 2020 06:56
 Operator : JC/SP
 Sample : L2018-04
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-2

Quant Time: Mar 30 05:47:50 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X032720W.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 27 09:35:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	1078628	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.84	114	1762703	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	1691417	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	915775	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.03	65	734258	48.27	ug/l	0.00
Spiked Amount			Recovery	=	96.54%	
35) Dibromofluoromethane	5.47	113	563734	50.03	ug/l	0.00
Spiked Amount			Recovery	=	100.06%	
50) Toluene-d8	8.70	98	2145080	51.71	ug/l	0.00
Spiked Amount			Recovery	=	103.42%	
62) 4-Bromofluorobenzene	11.12	95	933437	55.37	ug/l	0.00
Spiked Amount			Recovery	=	110.74%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.42	43	27812	4.255	ug/l	94
17) Carbon Disulfide	2.55	76	17465	0.562	ug/l	99
19) Methyl tert-butyl Ether	3.17	73	126728	3.209	ug/l	99
27) cis-1,2-Dichloroethene	4.58	96	8079	0.618	ug/l	96
65) Chlorobenzene	10.12	112	11293	0.335	ug/l	95

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

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