

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120221\
 Data File : VX025457.D
 Acq On : 03 Dec 2021 03:17
 Operator : JC/MD
 Sample : M4918-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-37

Quant Time: Dec 03 06:15:23 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 17 15:36:16 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.556	168	102786	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	174473	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	155262	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	68599	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	62793	57.96	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	115.92%
35) Dibromofluoromethane	5.391	113	54839	51.23	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.46%
50) Toluene-d8	8.653	98	200321	49.77	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.54%
62) 4-Bromofluorobenzene	11.079	95	68761	45.16	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.32%
Target Compounds						
						Qvalue
16) Acetone	2.380	43	1331	2.21	ug/l	97
17) Carbon Disulfide	2.514	76	3264	1.22	ug/l #	93
19) Methyl tert-butyl Ether	3.117	73	31412	9.46	ug/l	97
22) Diisopropyl ether	3.764	45	896	0.27	ug/l #	81
40) Benzene	6.038	78	1355	0.29	ug/l	91

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

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