

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061721\
 Data File : VX022739.D
 Acq On : 17 Jun 2021 14:41
 Operator : JC/MD
 Sample : M2750-18
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW178I-20210616

Quant Time: Jun 18 01:17:17 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 10 12:20:00 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	54086	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.769	114	111001	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	97075	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	37499	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	47404	48.733	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.460%
35) Dibromofluoromethane	5.397	113	34046	48.073	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.140%
50) Toluene-d8	8.653	98	136546	48.909	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.820%
62) 4-Bromofluorobenzene	11.085	95	46409	43.953	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.900%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	1256	2.884	ug/l	# 69
19) Methyl tert-butyl Ether	3.123	73	2531	1.111	ug/l	# 73
30) Chloroform	5.111	83	700	0.482	ug/l	# 82
44) Trichloroethene	7.129	130	1130	1.463	ug/l	99
64) Tetrachloroethene	9.275	164	316	0.515	ug/l	91

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

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