

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032521\
 Data File : VX021512.D
 Acq On : 26 Mar 2021 02:04
 Operator : JC/MD
 Sample : M1770-23
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-102I

Quant Time: Mar 26 06:34:13 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 24 12:25:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.623	168	57950	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.829	114	107648	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	108864	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	47526	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.035	65	51634	53.55	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 107.10%		
35) Dibromofluoromethane	5.470	113	39094	51.13	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 102.26%		
50) Toluene-d8	8.694	98	136356	50.04	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 100.08%		
62) 4-Bromofluorobenzene	11.118	95	53416	50.42	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 100.84%		
Target Compounds						
						Qvalue
16) Acetone	2.423	43	1820	4.99	ug/l	91
19) Methyl tert-butyl Ether	3.170	73	3524	1.62	ug/l	92
27) cis-1,2-Dichloroethene	4.564	96	5698	7.65	ug/l	99
30) Chloroform	5.176	83	1040	0.79	ug/l #	76
44) Trichloroethene	7.188	130	22072	27.93	ug/l	93
64) Tetrachloroethene	9.318	164	39703	54.44	ug/l	93

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

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