

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
 Data File : VX021598.D
 Acq On : 30 Mar 2021 00:17
 Operator : JC/MD
 Sample : M1843-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW178I-20210324

Quant Time: Mar 30 05:21:57 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	5.629	168	58844	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.829	114	109033	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	107842	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	44026	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.029	65	50930	52.02	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.04%
35) Dibromofluoromethane	5.464	113	39531	51.04	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.08%
50) Toluene-d8	8.694	98	136335	49.40	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.80%
62) 4-Bromofluorobenzene	11.118	95	51880	48.35	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.70%
Target Compounds						
					Qvalue	
16) Acetone	2.423	43	1380	3.73	ug/l #	82
19) Methyl tert-butyl Ether	3.164	73	2137	0.96	ug/l #	87
30) Chloroform	5.176	83	601	0.45	ug/l	94
44) Trichloroethene	7.188	130	906	1.13	ug/l	98
64) Tetrachloroethene	9.318	164	382	0.53	ug/l #	76

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

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