

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031822\
 Data File : VX027541.D
 Acq On : 18 Mar 2022 20:26
 Operator : JC/MD
 Sample : N1998-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-138

Quant Time: Mar 19 03:10:23 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 17 15:02:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	320440	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	531809	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	465487	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	192512	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	194661	49.312	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	98.620%
35) Dibromofluoromethane	5.391	113	177183	51.784	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	103.560%
50) Toluene-d8	8.653	98	620592	47.931	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	95.860%
62) 4-Bromofluorobenzene	11.085	95	208250	43.760	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	87.520%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	5592	3.475	ug/l	97
17) Carbon Disulfide	2.514	76	1936	0.256	ug/l #	89
19) Methyl tert-butyl Ether	3.117	73	1523441	151.750	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031822\
Data File : VX027541.D
Acq On : 18 Mar 2022 20:26
Operator : JC/MD
Sample : N1998-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-138

Quant Time: Mar 19 03:10:23 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
Quant Title : SW846 8260
QLast Update : Thu Mar 17 15:02:59 2022
Response via : Initial Calibration

