

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031822\
 Data File : VX027522.D
 Acq On : 18 Mar 2022 12:57
 Operator : JC/MD
 Sample : N1997-17DL 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-121DL

Quant Time: Mar 19 03:06:13 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	337633	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	559910	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	514390	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	245740	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	208070	50.025	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	100.060%
35) Dibromofluoromethane	5.391	113	182545	50.674	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	101.340%
50) Toluene-d8	8.653	98	657165	48.208	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.420%
62) 4-Bromofluorobenzene	11.085	95	242929	48.485	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	96.960%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.965	59	55360	69.839	ug/l	99
16) Acetone	2.380	43	2833	1.671	ug/l	90
17) Carbon Disulfide	2.514	76	3810	0.478	ug/l	98
19) Methyl tert-butyl Ether	3.117	73	361712	34.195	ug/l	99
40) Benzene	6.043	78	3185	0.226	ug/l #	84

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

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