

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031622\
 Data File : VN071331.D
 Acq On : 16 Mar 2022 19:37
 Operator : JC\MD
 Sample : N1957-13
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 D-13

Quant Time: Mar 17 05:21:02 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 08 06:52:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	886846	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.969	114	1469013	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1375731	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	513625	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	560073	46.079	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	92.160%
35) Dibromofluoromethane	8.022	113	420367	45.347	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	90.700%
50) Toluene-d8	10.439	98	1790241	47.965	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	95.940%
62) 4-Bromofluorobenzene	12.727	95	673856	53.317	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	106.640%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	13188	2.230	ug/l	99
18) Methyl Acetate	4.857	43	17703	1.574	ug/l	93
64) Tetrachloroethene	10.980	164	13107	1.366	ug/l	89

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

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