

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031822\
 Data File : VN071379.D
 Acq On : 18 Mar 2022 17:55
 Operator : JC\MD
 Sample : N2009-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31722

Quant Time: Mar 19 07:10:10 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	8.087	168	815254	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.969	114	1369984	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1306695	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.669	152	526360	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	527030	47.168	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	94.340%
35) Dibromofluoromethane	8.016	113	405454	46.900	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	93.800%
50) Toluene-d8	10.439	98	1690327	48.562	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	97.120%
62) 4-Bromofluorobenzene	12.727	95	667341	56.618	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	113.240%
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	127850	23.517	ug/l #	89
17) Carbon Disulfide	4.540	76	103720	4.458	ug/l	97
18) Methyl Acetate	4.851	43	234717	16.394	ug/l	96
25) 2-Butanone	7.334	43	151329	17.279	ug/l	95
51) 4-Methyl-2-Pentanone	10.328	43	18719	1.058	ug/l	94
70) Styrene	12.292	104	21071	0.748	ug/l #	88
95) Naphthalene	15.504	128	12253	5.524	ug/l	97

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

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