

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031622\
 Data File : VN071322.D
 Acq On : 16 Mar 2022 16:00
 Operator : JC\MD
 Sample : N1886-03 10X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 30969

Quant Time: Mar 17 05:18:42 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	8.081	168	807024	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1326603	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1253500	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	493630	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	504037	45.571	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	91.140%
35) Dibromofluoromethane	8.022	113	387644	46.306	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	92.620%
50) Toluene-d8	10.439	98	1654691	49.093	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	98.180%
62) 4-Bromofluorobenzene	12.727	95	627765	55.002	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	110.000%
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	151454	28.142	ug/l	99
18) Methyl Acetate	4.857	43	41752	3.332	ug/l	99
25) 2-Butanone	7.345	43	16915	1.951	ug/l #	86
51) 4-Methyl-2-Pentanone	10.328	43	190825	11.139	ug/l	99
59) 2-Hexanone	11.080	43	10228	0.875	ug/l	88
68) m/p-Xylenes	11.945	106	44114	2.533	ug/l	95

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

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