

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031622\
 Data File : VN071324.D
 Acq On : 16 Mar 2022 16:48
 Operator : JC\MD
 Sample : N1886-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 2822-A

Quant Time: Mar 17 05:19:12 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	934074	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1546797	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1469952	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	586432	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	596071	46.561	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	93.120%		
35) Dibromofluoromethane	8.022	113	454401	46.554	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	93.100%		
50) Toluene-d8	10.439	98	1924323	48.965	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	97.920%		
62) 4-Bromofluorobenzene	12.727	95	717935	53.948	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	107.900%		
Target Compounds						Qvalue
11) Tert butyl alcohol	5.357	59	33537	13.335	ug/l #	90
16) Acetone	4.281	43	1851669	297.268	ug/l	98
17) Carbon Disulfide	4.528	76	37227	1.396	ug/l	97
18) Methyl Acetate	4.851	43	217566	13.353	ug/l	96
20) Methylene Chloride	5.093	84	10546	1.301	ug/l #	88
25) 2-Butanone	7.334	43	216162	21.543	ug/l	95
29) Tetrahydrofuran	7.692	42	17916	2.577	ug/l	90
30) Chloroform	7.816	83	8648	0.428	ug/l	96
51) 4-Methyl-2-Pentanone	10.328	43	271272	13.581	ug/l	100
52) Toluene	10.504	92	13954	0.481	ug/l	93
59) 2-Hexanone	11.080	43	52845	3.877	ug/l	98
68) m/p-Xylenes	11.945	106	84667	4.146	ug/l	97

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

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