

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020122\
 Data File : VN070882.D
 Acq On : 1 Feb 2022 19:53
 Operator : JC/MD
 Sample : N1342-17
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31536

Quant Time: Feb 02 00:38:51 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 18 13:32:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1397585	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.966	114	2121373	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.741	117	1894311	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.672	152	863590	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	799521	39.917	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	79.840%		
35) Dibromofluoromethane	8.021	113	602761	45.806	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	91.620%		
50) Toluene-d8	10.441	98	2566072	46.290	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	92.580%		
62) 4-Bromofluorobenzene	12.728	95	941770	47.675	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	95.340%		
Target Compounds						
					Qvalue	
16) Acetone	4.291	43	136654	9.830	ug/l	96
20) Methylene Chloride	5.095	84	19349	1.057	ug/l	90
40) Benzene	8.459	78	324647	4.730	ug/l	99
43) Isopropyl Acetate	8.569	43	295165	6.571	ug/l #	77
51) 4-Methyl-2-Pentanone	10.331	43	329423	11.493	ug/l	95
52) Toluene	10.505	92	371628	9.051	ug/l	97
67) Ethyl Benzene	11.843	91	42625	0.560	ug/l	98
68) m/p-Xylenes	11.948	106	100691	3.637	ug/l	92
69) o-Xylene	12.278	106	28880	1.051	ug/l	93
84) 1,2,4-Trimethylbenzene	13.367	105	69427	1.093	ug/l	86
95) Naphthalene	15.499	128	42063	5.795	ug/l #	91

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

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