

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121521\
 Data File : VN070121.D
 Acq On : 15 Dec 2021 19:17
 Operator : JC/MD
 Sample : M4970-13RE
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SW-2RE

Quant Time: Dec 16 05:41:20 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	1110460	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.968	114	2145477	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2502073	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1056837	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	888563	51.825	ug/l	0.00
Spiked Amount	50.000	Range 61 - 141	Recovery	= 103.640%		
35) Dibromofluoromethane	8.024	113	621743	46.265	ug/l	0.00
Spiked Amount	50.000	Range 69 - 133	Recovery	= 92.520%		
50) Toluene-d8	10.443	98	2864903	53.092	ug/l	0.00
Spiked Amount	50.000	Range 65 - 126	Recovery	= 106.180%		
62) 4-Bromofluorobenzene	12.731	95	1246967	63.787	ug/l	0.00
Spiked Amount	50.000	Range 58 - 135	Recovery	= 127.580%		
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	299653	29.054	ug/l	96
25) 2-Butanone	7.351	43	64225	4.583	ug/l	94
40) Benzene	8.464	78	198477	3.179	ug/l	99
51) 4-Methyl-2-Pentanone	10.336	43	32336	1.108	ug/l	95
52) Toluene	10.505	92	26239	0.688	ug/l	95
67) Ethyl Benzene	11.846	91	20642	0.224	ug/l	94
78) n-propylbenzene	12.921	91	27541	0.263	ug/l	89
86) p-Isopropyltoluene	13.616	119	76192	1.072	ug/l	# 77
95) Naphthalene	15.509	128	12512	4.208	ug/l	99

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

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