

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121721\
 Data File : VN070170.D
 Acq On : 17 Dec 2021 17:52
 Operator : JC/MD
 Sample : M5039-06
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-25

Quant Time: Dec 18 05:47:04 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	864537	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1453921	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	1308555	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	454866	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	709632	53.162	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 106.320%	
35) Dibromofluoromethane	8.024	113	460847	50.603	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 101.200%	
50) Toluene-d8	10.443	98	1785742	48.834	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 97.660%	
62) 4-Bromofluorobenzene	12.734	95	587193	44.324	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 88.640%	
Target Compounds						
					Qvalue	
3) Chloromethane	2.301	50	3044	0.241	ug/l	97
11) Tert butyl alcohol	5.388	59	758808	300.734	ug/l	100
16) Acetone	4.299	43	35936	4.475	ug/l	100
19) Methyl tert-butyl Ether	5.621	73	2120790	66.233	ug/l	96
22) Diisopropyl ether	6.503	45	150140	4.140	ug/l	90
67) Ethyl Benzene	11.846	91	28139	0.584	ug/l	94
68) m/p-Xylenes	11.950	106	29884	1.702	ug/l	91
69) o-Xylene	12.280	106	3877	0.219	ug/l	90
78) n-propylbenzene	12.921	91	16521	0.367	ug/l	97
80) 1,3,5-Trimethylbenzene	13.061	105	9353	0.287	ug/l	98
84) 1,2,4-Trimethylbenzene	13.369	105	39058	1.259	ug/l	99

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

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