

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122921\
 Data File : VN070382.D
 Acq On : 30 Dec 2021 7:09
 Operator : JC/MD
 Sample : M5007-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TOLL-MW-05RE

Quant Time: Dec 30 08:03:26 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
 Quant Title : SW846 8260
 QLast Update : Mon Dec 27 18:47:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	824620	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1449211	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	1270761	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	403592	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	661938	55.829	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	111.660%
35) Dibromofluoromethane	8.021	113	466696	53.224	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	106.440%
50) Toluene-d8	10.443	98	1754042	48.968	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	97.940%
62) 4-Bromofluorobenzene	12.733	95	591285	45.703	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	91.400%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	84315	10.955	ug/l	99
17) Carbon Disulfide	4.532	76	10232	0.446	ug/l	97
18) Methyl Acetate	4.864	43	10799	0.485	ug/l #	64
19) Methyl tert-butyl Ether	5.621	73	72636	2.329	ug/l	95
25) 2-Butanone	7.351	43	14032	1.448	ug/l	91

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

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