

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120721\
 Data File : VN069824.D
 Acq On : 7 Dec 2021 19:14
 Operator : JC/MD
 Sample : M4956-23
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW178I-20211206

Quant Time: Dec 08 04:39:00 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	1338454	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2402246	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2553860	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	945533	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	1037263	50.192	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 100.380%	
35) Dibromofluoromethane	8.024	113	718099	47.723	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 95.440%	
50) Toluene-d8	10.440	98	3108625	51.451	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 102.900%	
62) 4-Bromofluorobenzene	12.731	95	1185122	54.143	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 108.280%	
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	19368	1.558	ug/l	92
19) Methyl tert-butyl Ether	5.624	73	23781	0.480	ug/l	96
30) Chloroform	7.823	83	8254	0.281	ug/l #	83
44) Trichloroethene	9.215	130	11176	0.670	ug/l	82
64) Tetrachloroethene	10.982	164	4963	0.287	ug/l	90

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

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