

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120821\
 Data File : VN069869.D
 Acq On : 8 Dec 2021 18:36
 Operator : JC/MD
 Sample : M4957-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW179D-20211206

Quant Time: Dec 09 11:44:35 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	1282247	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2306245	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2501139	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	969198	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	1000017	50.511	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	101.020%
35) Dibromofluoromethane	8.024	113	691324	47.856	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	95.720%
50) Toluene-d8	10.440	98	3007055	51.842	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	103.680%
62) 4-Bromofluorobenzene	12.731	95	1191466	56.699	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	113.400%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	22101	1.856	ug/l	91
19) Methyl tert-butyl Ether	5.613	73	18881	0.398	ug/l	96
27) cis-1,2-Dichloroethene	7.329	96	254706	15.891	ug/l	96
44) Trichloroethene	9.217	130	1092279	68.195	ug/l	92
64) Tetrachloroethene	10.977	164	152140	8.971	ug/l	96

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

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