

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
 Data File : VN069558.D
 Acq On : 18 Nov 2021 17:25
 Operator : JC/MD
 Sample : M4739-05
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-23A

Quant Time: Nov 19 03:19:41 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 09 18:18:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	1366934	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	2374098	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	2382935	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	890291	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	1017364	51.331	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	102.660%
35) Dibromofluoromethane	8.024	113	695472	48.843	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	97.680%
50) Toluene-d8	10.443	98	3035179	54.199	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	108.400%
62) 4-Bromofluorobenzene	12.733	95	1083631	52.673	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	105.340%
Target Compounds						
						Qvalue
16) Acetone	4.301	43	16627	1.555	ug/l	# 86
19) Methyl tert-butyl Ether	5.621	73	9946161	212.130	ug/l	96
22) Diisopropyl ether	6.498	45	240667	4.561	ug/l	91
42) 1,2-Dichloroethane	8.531	62	20039	0.830	ug/l	84

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

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