

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
 Data File : VN069555.D
 Acq On : 18 Nov 2021 16:09
 Operator : JC/MD
 Sample : M4739-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-6R

Quant Time: Nov 19 03:19:06 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	1369317	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	2378920	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	2365178	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	851567	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	1012288	50.986	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	101.980%
35) Dibromofluoromethane	8.024	113	702141	49.212	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	98.420%
50) Toluene-d8	10.443	98	3039742	54.170	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	108.340%
62) 4-Bromofluorobenzene	12.734	95	1053584	51.109	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	102.220%
Target Compounds						
						Qvalue
16) Acetone	4.301	43	59890	5.591	ug/l	92
19) Methyl tert-butyl Ether	5.624	73	1407520	29.967	ug/l	96
22) Diisopropyl ether	6.503	45	64365	1.218	ug/l	90
40) Benzene	8.464	78	12429	0.201	ug/l	91

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

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