

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
 Data File : VN069559.D
 Acq On : 18 Nov 2021 17:51
 Operator : JC/MD
 Sample : M4739-06
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-33A

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/19/2021
 Supervised By : Semsettin Yesilyurt 11/19/2021

Quant Time: Nov 19 03:19:51 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	1370817	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	2410233	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	2373261	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	858546	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	1022886	51.463	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 102.920%			
35) Dibromofluoromethane	8.024	113	700799	48.480	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 96.960%			
50) Toluene-d8	10.443	98	3052232	53.686	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 107.380%			
62) 4-Bromofluorobenzene	12.733	95	1039923	49.791	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 99.580%			
Target Compounds						
16) Acetone	4.301	43	20532	1.915	ug/l	99
19) Methyl tert-butyl Ether	5.621	73	195096	4.149	ug/l	94
24) 1,1-Dichloroethane	6.383	63	19806m	0.735	ug/l	
27) cis-1,2-Dichloroethene	7.329	96	62931	4.091	ug/l	96

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

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