

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111921\
 Data File : VN069596.D
 Acq On : 19 Nov 2021 19:55
 Operator : JC/MD
 Sample : M4768-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-3

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/22/2021
 Supervised By : Mahesh Dadoda 11/26/2021

Quant Time: Nov 22 05:09:35 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	1452396	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2521207	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	2518048	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	999805	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	1080911	51.328	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 102.660%			
35) Dibromofluoromethane	8.024	113	746665	49.379	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 98.760%			
50) Toluene-d8	10.443	98	3184634	53.549	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 107.100%			
62) 4-Bromofluorobenzene	12.731	95	1185008	54.240	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 108.480%			
Target Compounds						
16) Acetone	4.293	43	15393m	1.355	ug/l	
52) Toluene	10.508	92	99299	2.434	ug/l	97
67) Ethyl Benzene	11.846	91	68856	0.797	ug/l	98
68) m/p-Xylenes	11.950	106	123668	3.890	ug/l	93
69) o-Xylene	12.280	106	54516	1.720	ug/l	91
80) 1,3,5-Trimethylbenzene	13.061	105	16920	0.263	ug/l	99
83) tert-Butylbenzene	13.321	119	19775	0.351	ug/l	83
84) 1,2,4-Trimethylbenzene	13.369	105	60515	0.974	ug/l	99
85) sec-Butylbenzene	13.501	105	50451	0.640	ug/l #	51

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

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