

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120321\
 Data File : VN069744.D
 Acq On : 4 Dec 2021 3:06
 Operator : JC/MD
 Sample : M4919-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FDR-MW-3-20211202

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/04/2021
 Supervised By : Mahesh Dadoda 12/06/2021

Quant Time: Dec 04 05:41:30 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	1524853	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2683372	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2759971	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1055090	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	1184231	50.30	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 100.60%			
35) Dibromofluoromethane	8.024	113	812244	48.32	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 96.64%			
50) Toluene-d8	10.440	98	3446148	51.06	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 102.12%			
62) 4-Bromofluorobenzene	12.731	95	1269376	51.92	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 103.84%			
Target Compounds						Qvalue
17) Carbon Disulfide	4.543	76	30475m	0.81	ug/l	93
19) Methyl tert-butyl Ether	5.618	73	15876m	0.28	ug/l	
83) tert-Butylbenzene	13.324	119	226634	3.39	ug/l	

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

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