

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122121\
 Data File : VN070272.D
 Acq On : 21 Dec 2021 17:09
 Operator : JC/MD
 Sample : M5039-11
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-31

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/22/2021
 Supervised By : Semsettin Yesilyurt 12/23/2021

Quant Time: Dec 21 20:11:51 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	821277	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1370610	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1186420	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	391933	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	691282	54.515	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 109.040%			
35) Dibromofluoromethane	8.021	113	454992	52.997	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 106.000%			
50) Toluene-d8	10.440	98	1727439	50.111	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 100.220%			
62) 4-Bromofluorobenzene	12.731	95	553187	44.295	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 88.600%			
Target Compounds						
3) Chloromethane	2.306	50	3968	0.331	ug/l	94
16) Acetone	4.293	43	30375	3.982	ug/l	93
64) Tetrachloroethene	10.971	164	3802m	0.473	ug/l	
68) m/p-Xylenes	11.956	106	6233	0.392	ug/l	95
84) 1,2,4-Trimethylbenzene	13.369	105	12458	0.466	ug/l	97

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

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