

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120424\
 Data File : VN085096.D
 Acq On : 04 Dec 2024 16:01
 Operator : JC\MD
 Sample : P5072-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-738

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2024
 Supervised By : Mahesh Dadoda 12/05/2024

Quant Time: Dec 04 22:08:55 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112224W.M
 Quant Title : SW846 8260
 QLast Update : Sat Nov 23 02:54:10 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.224	168	153255	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	283554	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	254079	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	97320	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	113670	52.070	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 104.140%			
35) Dibromofluoromethane	8.165	113	94502	49.421	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 98.840%			
50) Toluene-d8	10.565	98	333904	48.725	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 97.440%			
62) 4-Bromofluorobenzene	12.847	95	116086	45.298	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 90.600%			
Target Compounds						
16) Acetone	4.430	43	37268	58.618	ug/l	96
43) Isopropyl Acetate	8.688	43	150458m	30.762	ug/l	
51) 4-Methyl-2-Pentanone	10.441	43	10106m	4.524	ug/l	

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

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