

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120424\
 Data File : VN085095.D
 Acq On : 04 Dec 2024 15:37
 Operator : JC\MD
 Sample : P5072-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MH-737

Manual Integrations
 APPROVED

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

Quant Time: Dec 04 22:08:26 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	162661	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	293085	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	264444	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	100644	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	116365	50.223	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 100.440%			
35) Dibromofluoromethane	8.165	113	95813	48.477	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 96.960%			
50) Toluene-d8	10.565	98	348964	49.267	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 98.540%			
62) 4-Bromofluorobenzene	12.847	95	119216	45.006	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 90.020%			
Target Compounds						
16) Acetone	4.424	43	36153	53.576	ug/l	Qvalue 92
43) Isopropyl Acetate	8.688	43	141433m	27.977	ug/l	
51) 4-Methyl-2-Pentanone	10.447	43	10191	4.413	ug/l #	71

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

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