

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102824\
 Data File : VN084548.D
 Acq On : 28 Oct 2024 18:41
 Operator : JC\MD
 Sample : P4567-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-1

Manual Integrations
 APPROVED

Reviewed By :Mahesh Dadoda 10/29/2024
 Supervised By :Semsettin Yesilyurt 10/29/2024

Quant Time: Oct 29 01:32:58 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N093024W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 01 07:11:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	166485	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	302607	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	275908	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	110514	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	124189	50.311	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 100.620%			
35) Dibromofluoromethane	8.165	113	99359	49.805	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 99.600%			
50) Toluene-d8	10.565	98	374968	51.087	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 102.180%			
62) 4-Bromofluorobenzene	12.847	95	128892	48.203	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 96.400%			
Target Compounds					Qvalue	
16) Acetone	4.430	43	35623	33.594	ug/l	99
20) Methylene Chloride	5.289	84	3254	1.510	ug/l #	82
43) Isopropyl Acetate	8.688	43	146021m	25.077	ug/l	

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

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