

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102424\
 Data File : VN084496.D
 Acq On : 24 Oct 2024 18:21
 Operator : JC\MD
 Sample : P4468-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ETGI-329

Manual Integrations
 APPROVED

Reviewed By :Mahesh Dadoda 10/25/2024
 Supervised By :Semsettin Yesilyurt 10/25/2024

Quant Time: Oct 25 02:05:27 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	185646	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	334594	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	287090	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	108396	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	130241	47.317	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	94.640%		
35) Dibromofluoromethane	8.165	113	105453	47.806	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.620%		
50) Toluene-d8	10.565	98	397697	49.004	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	98.000%		
62) 4-Bromofluorobenzene	12.847	95	128382	43.422	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	86.840%		
Target Compounds					Qvalue	
16) Acetone	4.436	43	27199	23.003	ug/l	93
20) Methylene Chloride	5.271	84	4181	1.739	ug/l	94
43) Isopropyl Acetate	8.688	43	247593m	38.455	ug/l	

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

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