

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102424\
 Data File : VN084497.D
 Acq On : 24 Oct 2024 18:45
 Operator : JC\MD
 Sample : P4468-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ETGI-345

Manual Integrations
 APPROVED

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

Quant Time: Oct 25 02:05:51 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	182228	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	332892	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	290809	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	111115	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	128444	47.539	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	95.080%		
35) Dibromofluoromethane	8.171	113	102681	46.787	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	93.580%		
50) Toluene-d8	10.565	98	401136	49.681	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	99.360%		
62) 4-Bromofluorobenzene	12.847	95	132535	45.056	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	90.120%		
Target Compounds						
16) Acetone	4.436	43	26217	22.588	ug/l	99
20) Methylene Chloride	5.283	84	4640	1.967	ug/l	89
43) Isopropyl Acetate	8.688	43	236004m	36.843	ug/l	

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

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