

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081324\
 Data File : VN083287.D
 Acq On : 14 Aug 2024 08:52
 Operator : JC\MD
 Sample : P3466-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 924-K1-WP0-0.25-080224

Manual Integrations
 APPROVED

Reviewed By : John Carlone 08/14/2024
 Supervised By : Mahesh Dadoda 08/14/2024

Quant Time: Aug 14 09:34:00 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080724W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 08 06:30:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	128494	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	251972	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	264206	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	114718	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	108106	59.108	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 118.220%			
35) Dibromofluoromethane	8.165	113	86756	55.162	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 110.320%			
50) Toluene-d8	10.565	98	316687	53.980	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 107.960%			
62) 4-Bromofluorobenzene	12.847	95	132070	57.743	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 115.480%			
Target Compounds						
16) Acetone	4.447	43	3340m	4.667	ug/l	
29) Tetrahydrofuran	7.847	42	5049	7.108	ug/l #	86
52) Toluene	10.629	92	2504	0.559	ug/l	93

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

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