

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091824\
 Data File : VN084021.D
 Acq On : 19 Sep 2024 08:56
 Operator : JC\MD
 Sample : P4097-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RT2299

Manual Integrations
 APPROVED

Reviewed By : John Carlone 09/19/2024
 Supervised By : Mahesh Dadoda 09/19/2024

Quant Time: Sep 19 09:18:03 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 11 15:18:56 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	154005	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	273890	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	236306	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	99862	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	121685	52.401	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 104.800%			
35) Dibromofluoromethane	8.165	113	87387	49.838	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 99.680%			
50) Toluene-d8	10.565	98	331329	54.660	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 109.320%			
62) 4-Bromofluorobenzene	12.847	95	120800	48.303	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 96.600%			
Target Compounds						
16) Acetone	4.430	43	18868	20.268	ug/l	94
20) Methylene Chloride	5.277	84	5767	3.027	ug/l	95
25) 2-Butanone	7.483	43	9016	7.369	ug/l	96
43) Isopropyl Acetate	8.688	43	224900m	51.808	ug/l	

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

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