

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091824\
 Data File : VN084016.D
 Acq On : 19 Sep 2024 06:57
 Operator : JC\MD
 Sample : P4087-32
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 NB-614-COMP-06

Manual Integrations
 APPROVED

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

Quant Time: Sep 19 07:38:35 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	167810	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	289798	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	249571	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	100841	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	130089	51.412	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 102.820%			
35) Dibromofluoromethane	8.165	113	89393	48.184	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 96.360%			
50) Toluene-d8	10.565	98	347218	54.137	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 108.280%			
62) 4-Bromofluorobenzene	12.847	95	128067	48.398	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 96.800%			
Target Compounds						
16) Acetone	4.430	43	16499	16.266	ug/l	91
20) Methylene Chloride	5.271	84	5320	2.563	ug/l #	85
25) 2-Butanone	7.488	43	9157	6.868	ug/l	99
43) Isopropyl Acetate	8.688	43	243021m	52.909	ug/l	

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

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