

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091824\
 Data File : VN084019.D
 Acq On : 19 Sep 2024 08:09
 Operator : JC\MD
 Sample : P4088-14
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 NB-614-COMP-09

Manual Integrations
 APPROVED

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

Quant Time: Sep 19 08:34:54 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	167404	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	290486	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	248966	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	103847	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.576	65	129529	51.315	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	102.620%	
35) Dibromofluoromethane	8.165	113	91113	48.995	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	97.980%	
50) Toluene-d8	10.565	98	347605	54.069	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	108.140%	
62) 4-Bromofluorobenzene	12.847	95	125432	47.290	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	=	94.580%	
Target Compounds						
16) Acetone	4.436	43	31553	31.182	ug/l	93
20) Methylene Chloride	5.283	84	5475	2.644	ug/l #	85
25) 2-Butanone	7.488	43	14931	11.227	ug/l	94
43) Isopropyl Acetate	8.688	43	234140m	50.855	ug/l	

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

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