

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091024\
 Data File : VN083761.D
 Acq On : 10 Sep 2024 18:56
 Operator : JC\MD
 Sample : P3907-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PL-701-COMP-54

Quant Time: Sep 11 03:18:21 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 11 02:47:22 2024
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :John Carlone 09/11/2024
 Supervised By :Semsettin Yesilyurt 09/11/2024

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	114204	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	241358	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	269513	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	129958	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	92638	53.796	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 107.600%			
35) Dibromofluoromethane	8.165	113	72623	47.001	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 94.000%			
50) Toluene-d8	10.565	98	269164	50.390	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 100.780%			
62) 4-Bromofluorobenzene	12.847	95	126201	57.264	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 114.520%			
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	41679	60.376	ug/l	100
20) Methylene Chloride	5.259	84	6258	4.430	ug/l #	90
37) Ethyl Acetate	7.559	43	3084	1.461	ug/l #	87
43) Isopropyl Acetate	8.689	43	197149m	51.536	ug/l	

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

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