

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091024\
 Data File : VN083786.D
 Acq On : 11 Sep 2024 10:50
 Operator : JC\MD
 Sample : P3906-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SU-701-COMP-02

Quant Time: Sep 11 11:08:44 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 11 03:05:46 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	117297	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	253147	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	282340	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	134617	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	99207	56.091	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 112.180%			
35) Dibromofluoromethane	8.165	113	76610	47.272	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 94.540%			
50) Toluene-d8	10.565	98	283492	50.600	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 101.200%			
62) 4-Bromofluorobenzene	12.847	95	133927	57.940	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 115.880%			
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	38982	54.980	ug/l	93
18) Methyl Acetate	5.030	43	9864	5.007	ug/l	95
20) Methylene Chloride	5.259	84	6232	4.295	ug/l	87
25) 2-Butanone	7.489	43	4416	4.739	ug/l	90
43) Isopropyl Acetate	8.688	43	238567	59.459	ug/l #	75

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

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