

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091024\
 Data File : VN083785.D
 Acq On : 11 Sep 2024 10:26
 Operator : JC\MD
 Sample : P3906-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SU-701-COMP-01

Quant Time: Sep 11 10:44:45 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	8.218	168	121052	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	260253	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	292751	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	140235	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	101880	55.816	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	111.640%
35) Dibromofluoromethane	8.165	113	81700	49.036	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.080%
50) Toluene-d8	10.565	98	297664	51.679	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.360%
62) 4-Bromofluorobenzene	12.847	95	135686	57.098	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	114.200%
Target Compounds						
						Qvalue
16) Acetone	4.430	43	39602	54.122	ug/l	97
18) Methyl Acetate	5.030	43	9420	4.490	ug/l	93
20) Methylene Chloride	5.259	84	5726	3.824	ug/l	94
25) 2-Butanone	7.494	43	4605	4.788	ug/l #	83
43) Isopropyl Acetate	8.688	43	132388	32.095	ug/l #	90

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

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