

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091024\
 Data File : VN083784.D
 Acq On : 11 Sep 2024 10:02
 Operator : JC\MD
 Sample : P3892-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ARS20-005

Manual Integrations
 APPROVED

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

Quant Time: Sep 11 10:24:28 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	119507	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	260005	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	292488	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	142853	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	102024	56.617	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 113.240%			
35) Dibromofluoromethane	8.165	113	80831	48.561	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 97.120%			
50) Toluene-d8	10.565	98	302044	52.490	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 104.980%			
62) 4-Bromofluorobenzene	12.847	95	138265	58.239	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 116.480%			
Target Compounds						
16) Acetone	4.436	43	43095	59.657	ug/l	99
18) Methyl Acetate	5.036	43	10197m	5.109	ug/l	
20) Methylene Chloride	5.277	84	5568	3.767	ug/l	94
25) 2-Butanone	7.488	43	4188	4.411	ug/l	93
43) Isopropyl Acetate	8.688	43	272712	66.176	ug/l #	72

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

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