

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082422\
 Data File : VN074094.D
 Acq On : 24 Aug 2022 16:15
 Operator : JC\MD
 Sample : N4244-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-103S

Quant Time: Aug 25 07:54:51 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081522W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 16 00:45:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	571943	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	868502	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	729122	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	324007	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	291132	52.261	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 104.520%	
35) Dibromofluoromethane	7.951	113	262144	51.038	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.080%	
50) Toluene-d8	10.380	98	969376	48.731	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 97.460%	
62) 4-Bromofluorobenzene	12.674	95	306882	46.484	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 92.960%	
Target Compounds						
					Qvalue	
3) Chloromethane	2.269	50	1304	0.259	ug/l #	83
16) Acetone	4.240	43	11379	4.330	ug/l #	86
17) Carbon Disulfide	4.457	76	5419	0.505	ug/l #	88
40) Benzene	8.386	78	27438	1.219	ug/l	96
68) m/p-Xylenes	11.892	106	4509	0.431	ug/l	96

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

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