

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
 Data File : VN074277.D
 Acq On : 02 Sep 2022 06:09
 Operator : JC\MD
 Sample : N4354-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DUP082222

Quant Time: Sep 02 08:40:33 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N083022W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 31 07:02:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	190891	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	374733	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	364650	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	188071	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	150003	50.656	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 101.320%	
35) Dibromofluoromethane	7.951	113	124491	51.053	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.100%	
50) Toluene-d8	10.380	98	448735	46.729	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 93.460%	
62) 4-Bromofluorobenzene	12.669	95	174845	59.872	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 119.740%	
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	7362	3.721	ug/l	95
40) Benzene	8.386	78	8897	0.716	ug/l	99
67) Ethyl Benzene	11.786	91	8378	0.569	ug/l #	77
68) m/p-Xylenes	11.892	106	3926	0.718	ug/l	91
69) o-Xylene	12.227	106	1576	0.289	ug/l	83
84) 1,2,4-Trimethylbenzene	13.304	105	5041	0.399	ug/l	83
95) Naphthalene	15.433	128	19262	2.470	ug/l #	95

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

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