

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN011022\
 Data File : VN070554.D
 Acq On : 10 Jan 2022 19:51
 Operator : JC/MD
 Sample : N1051-07
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TW-1

Quant Time: Jan 11 06:07:14 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
 Quant Title : SW846 8260
 QLast Update : Mon Dec 27 18:47:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	978893	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1665735	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1483837	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.672	152	555029	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	735987	52.292	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 104.580%	
35) Dibromofluoromethane	8.021	113	521123	51.706	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 103.420%	
50) Toluene-d8	10.440	98	2129820	51.730	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 103.460%	
62) 4-Bromofluorobenzene	12.731	95	798491	53.697	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 107.400%	
Target Compounds						
					Qvalue	
13) Acrolein	4.044	56	8698	4.574	ug/l	97
16) Acetone	4.291	43	198071	21.679	ug/l	94
17) Carbon Disulfide	4.532	76	18357	0.675	ug/l	99
18) Methyl Acetate	4.865	43	25168	0.952	ug/l	93

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

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