

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120921\
 Data File : VN069913.D
 Acq On : 9 Dec 2021 15:55
 Operator : JC/MD
 Sample : M4940-07
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-23

Quant Time: Dec 10 06:16:40 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	1246329	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2338637	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2599926	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1066623	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	985362	51.205	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	102.420%
35) Dibromofluoromethane	8.024	113	719005	49.083	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	98.160%
50) Toluene-d8	10.440	98	3017750	51.305	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	102.620%
62) 4-Bromofluorobenzene	12.731	95	1253279	58.814	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	117.620%
Target Compounds						
						Qvalue
16) Acetone	4.264	43	50228	4.339	ug/l	# 60
17) Carbon Disulfide	4.540	76	6305	0.206	ug/l	100
40) Benzene	8.464	78	65712	0.966	ug/l	94
78) n-propylbenzene	12.921	91	33174m	0.314	ug/l	

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

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