

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121021\
 Data File : VN069974.D
 Acq On : 10 Dec 2021 21:56
 Operator : JC/MD
 Sample : M4970-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-22

Quant Time: Dec 11 05:26:23 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	1079295	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2077331	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	2370280	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	951829	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	862983	51.786	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 103.580%	
35) Dibromofluoromethane	8.024	113	642345	49.366	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 98.740%	
50) Toluene-d8	10.443	98	2742167	52.484	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 104.960%	
62) 4-Bromofluorobenzene	12.733	95	1156527	61.101	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 122.200%	
Target Compounds						
						Qvalue
11) Tert butyl alcohol	5.377	59	238251	75.636	ug/l	99
16) Acetone	4.293	43	35306	3.522	ug/l	92
19) Methyl tert-butyl Ether	5.621	73	1954858	48.903	ug/l	95
22) Diisopropyl ether	6.503	45	95674	2.113	ug/l #	89
29) Tetrahydrofuran	7.702	42	13848	1.674	ug/l #	87
40) Benzene	8.464	78	73527	1.216	ug/l	97

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

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