

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120921\
 Data File : VN069946.D
 Acq On : 10 Dec 2021 6:45
 Operator : JC/MD
 Sample : M4966-20
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 P14-11

Quant Time: Dec 10 08:16:39 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	1073554	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2109491	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2412005	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	972688	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	887722	53.556	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	107.120%
35) Dibromofluoromethane	8.024	113	649764	49.175	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	98.340%
50) Toluene-d8	10.443	98	2797197	52.721	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	105.440%
62) 4-Bromofluorobenzene	12.731	95	1145646	59.604	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	119.200%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	96709	9.699	ug/l	99
18) Methyl Acetate	4.862	43	47868	1.590	ug/l #	90
20) Methylene Chloride	5.103	84	133192	10.048	ug/l	87
43) Isopropyl Acetate	8.566	43	199319	4.506	ug/l #	78

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

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