

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120921\
 Data File : VN069945.D
 Acq On : 10 Dec 2021 6:19
 Operator : JC/MD
 Sample : M4966-13
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 P10-11

Quant Time: Dec 10 08:16:16 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	1118583	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2171570	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	2479427	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	998550	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	914033	52.923	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	105.840%
35) Dibromofluoromethane	8.024	113	669298	49.205	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	98.400%
50) Toluene-d8	10.440	98	2886435	52.848	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	105.700%
62) 4-Bromofluorobenzene	12.731	95	1177869	59.528	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	119.060%
Target Compounds						
						Qvalue
16) Acetone	4.293	43	316404	30.455	ug/l	97
18) Methyl Acetate	4.859	43	47891	1.527	ug/l #	73
20) Methylene Chloride	5.095	84	138971	10.062	ug/l	87
43) Isopropyl Acetate	8.566	43	199509	4.382	ug/l #	77

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

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