

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062123\
 Data File : VN078337.D
 Acq On : 21 Jun 2023 21:11
 Operator : JC\MD
 Sample : 03337-20
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-11

Quant Time: Jun 22 04:56:32 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060623W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 07 05:05:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.236	168	442146	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	852152	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	806564	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	272823	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	323646	54.900	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 109.800%	
35) Dibromofluoromethane	8.177	113	294165	54.596	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 109.200%	
50) Toluene-d8	10.571	98	1026142	49.402	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 98.800%	
62) 4-Bromofluorobenzene	12.853	95	342409	53.026	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 106.060%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	88622	52.228	ug/l	95
18) Methyl Acetate	5.047	43	9070	1.161	ug/l #	73
20) Methylene Chloride	5.295	84	19467	3.472	ug/l #	91
25) 2-Butanone	7.494	43	26073	9.992	ug/l #	70
37) Ethyl Acetate	7.571	43	67585	11.194	ug/l	96
43) Isopropyl Acetate	8.694	43	534862	46.323	ug/l #	70
51) 4-Methyl-2-Pentanone	10.453	43	152793	25.129	ug/l	91

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

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