

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062123\
 Data File : VN078336.D
 Acq On : 21 Jun 2023 20:47
 Operator : JC\MD
 Sample : 03337-17
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-15

Quant Time: Jun 22 04:56:06 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	468812	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	906141	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	835670	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	266552	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	325472	52.070	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 104.140%	
35) Dibromofluoromethane	8.177	113	307722	53.709	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 107.420%	
50) Toluene-d8	10.571	98	1074790	48.661	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 97.320%	
62) 4-Bromofluorobenzene	12.853	95	340676	49.615	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 99.220%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	165146	91.790	ug/l	96
20) Methylene Chloride	5.283	84	13099	2.203	ug/l	90
25) 2-Butanone	7.494	43	45292	16.370	ug/l	93
37) Ethyl Acetate	7.571	43	70606	10.998	ug/l	96
43) Isopropyl Acetate	8.694	43	511516	41.661	ug/l #	72
51) 4-Methyl-2-Pentanone	10.453	43	324463	50.184	ug/l	90

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

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