

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062423\
 Data File : VN078393.D
 Acq On : 24 Jun 2023 14:37
 Operator : JC\MD
 Sample : 03393-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SF-9

Quant Time: Jun 26 01:50:00 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	8.230	168	426856	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	793391	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	727670	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	212125	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	295246	51.877	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 103.760%	
35) Dibromofluoromethane	8.177	113	277030	55.224	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 110.440%	
50) Toluene-d8	10.571	98	948608	49.051	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 98.100%	
62) 4-Bromofluorobenzene	12.853	95	291103	48.420	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 96.840%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	78094	47.672	ug/l	96
20) Methylene Chloride	5.289	84	29838	5.512	ug/l	94
25) 2-Butanone	7.500	43	20302	8.059	ug/l #	80
37) Ethyl Acetate	7.571	43	55534	9.879	ug/l	96
43) Isopropyl Acetate	8.694	43	270391	25.152	ug/l #	82
51) 4-Methyl-2-Pentanone	10.453	43	151295	26.726	ug/l	90

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

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