

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071823\
 Data File : VN078568.D
 Acq On : 18 Jul 2023 18:04
 Operator : JC\MD
 Sample : 03654-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 72-12016

Quant Time: Jul 19 01:54:28 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	398284	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	755943	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	719475	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	259253	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	267758	50.147	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 100.300%		
35) Dibromofluoromethane	8.177	113	268517	52.326	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 104.660%		
50) Toluene-d8	10.571	98	894377	46.136	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 92.280%		
62) 4-Bromofluorobenzene	12.853	95	311863	52.228	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 104.460%		
Target Compounds						
					Qvalue	
16) Acetone	4.453	43	67836	40.786	ug/l	96
20) Methylene Chloride	5.283	84	18871	2.801	ug/l	91
25) 2-Butanone	7.506	43	42461	16.218	ug/l #	72
37) Ethyl Acetate	7.571	43	12587	2.038	ug/l #	95
43) Isopropyl Acetate	8.700	43	421928	41.491	ug/l #	75
51) 4-Methyl-2-Pentanone	10.453	43	145849	24.112	ug/l	94

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

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