

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071823\
 Data File : VN078560.D
 Acq On : 18 Jul 2023 14:50
 Operator : JC\MD
 Sample : 03625-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RBR200018

Quant Time: Jul 19 01:50:32 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	402410	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	739667	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	720444	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	268308	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	271352	50.299	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 100.600%		
35) Dibromofluoromethane	8.171	113	275855	54.939	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 109.880%		
50) Toluene-d8	10.571	98	894809	47.174	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 94.340%		
62) 4-Bromofluorobenzene	12.853	95	316889	54.238	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 108.480%		
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	97084	57.772	ug/l	94
20) Methylene Chloride	5.294	84	17188	2.402	ug/l	92
25) 2-Butanone	7.494	43	55959	21.154	ug/l #	83
37) Ethyl Acetate	7.571	43	16385	2.711	ug/l #	91
43) Isopropyl Acetate	8.694	43	279660	28.106	ug/l #	78
51) 4-Methyl-2-Pentanone	10.447	43	229437	38.766	ug/l	92

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

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