

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071823\
 Data File : VN078561.D
 Acq On : 18 Jul 2023 15:14
 Operator : JC\MD
 Sample : 03646-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GATE-1

Quant Time: Jul 19 01:51:05 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	8.230	168	408126	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	766574	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	743424	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	276569	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	284899	52.071	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 104.140%	
35) Dibromofluoromethane	8.177	113	281540	54.103	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 108.200%	
50) Toluene-d8	10.571	98	912266	46.406	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 92.820%	
62) 4-Bromofluorobenzene	12.853	95	328640	54.275	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 108.540%	
Target Compounds						
					Qvalue	
16) Acetone	4.453	43	64490	37.839	ug/l	94
20) Methylene Chloride	5.283	84	18910	2.712	ug/l	96
25) 2-Butanone	7.506	43	40785	15.202	ug/l #	85
37) Ethyl Acetate	7.577	43	12427	1.984	ug/l	97
43) Isopropyl Acetate	8.700	43	301059	29.194	ug/l #	80
51) 4-Methyl-2-Pentanone	10.453	43	141239	23.026	ug/l	86

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

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