

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071823\
 Data File : VN078562.D
 Acq On : 18 Jul 2023 15:38
 Operator : JC\MD
 Sample : 03646-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 YARD-1

Quant Time: Jul 19 01:51:34 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	8.236	168	409975	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	775557	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	836304	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	264305	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	282907	51.474	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 102.940%	
35) Dibromofluoromethane	8.177	113	277564	52.721	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 105.440%	
50) Toluene-d8	10.571	98	921724	46.344	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 92.680%	
62) 4-Bromofluorobenzene	12.853	95	381609	62.292	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 124.580%#	
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	69892	40.824	ug/l	94
20) Methylene Chloride	5.289	84	18335	2.574	ug/l #	87
25) 2-Butanone	7.494	43	43398	16.103	ug/l #	80
37) Ethyl Acetate	7.577	43	14087	2.223	ug/l #	91
43) Isopropyl Acetate	8.700	43	351848	33.724	ug/l #	78
51) 4-Methyl-2-Pentanone	10.453	43	142998	23.043	ug/l	90

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

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