

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072723\
 Data File : VN078672.D
 Acq On : 27 Jul 2023 20:11
 Operator : JC\MD
 Sample : 03770-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 HOWE-1-LP

Quant Time: Jul 28 01:12:32 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	8.236	168	424229	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	780459	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	753343	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	296762	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	277542	48.801	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.600%
35) Dibromofluoromethane	8.177	113	276408	52.172	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	104.340%
50) Toluene-d8	10.577	98	955566	47.744	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.480%
62) 4-Bromofluorobenzene	12.853	95	327834	53.178	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	106.360%
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	104122	58.774	ug/l	98
20) Methylene Chloride	5.295	84	22089	3.202	ug/l	93
25) 2-Butanone	7.500	43	57827	20.736	ug/l	91
37) Ethyl Acetate	7.571	43	17694	2.775	ug/l #	89
43) Isopropyl Acetate	8.700	43	467810	44.558	ug/l #	73
51) 4-Methyl-2-Pentanone	10.453	43	249719	39.988	ug/l	91

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

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