

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072723\
 Data File : VN078673.D
 Acq On : 27 Jul 2023 20:35
 Operator : JC\MD
 Sample : 03770-11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 HOWE-2-LP

Quant Time: Jul 28 01:13:02 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	738938	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	1364049	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	1271598	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	466357	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	494069	49.874	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.740%
35) Dibromofluoromethane	8.177	113	476382	51.447	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.900%
50) Toluene-d8	10.577	98	1690811	48.337	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.680%
62) 4-Bromofluorobenzene	12.853	95	594855	55.209	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	110.420%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	201196	65.201	ug/l	94
20) Methylene Chloride	5.300	84	37561	3.097	ug/l #	93
25) 2-Butanone	7.500	43	113993	23.468	ug/l #	81
37) Ethyl Acetate	7.571	43	39221	3.519	ug/l	97
43) Isopropyl Acetate	8.700	43	879038	47.905	ug/l #	75
51) 4-Methyl-2-Pentanone	10.453	43	482998	44.253	ug/l	93

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

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