

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072523\
 Data File : VN078648.D
 Acq On : 25 Jul 2023 15:36
 Operator : JC\MD
 Sample : 03738-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 80001

Quant Time: Jul 26 04:47:56 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	449086	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	817554	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	784804	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	295621	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	283508	47.091	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	94.180%
35) Dibromofluoromethane	8.171	113	281323	50.691	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.380%
50) Toluene-d8	10.571	98	979284	46.709	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	93.420%
62) 4-Bromofluorobenzene	12.853	95	343677	53.219	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	106.440%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	92320	49.227	ug/l	99
20) Methylene Chloride	5.294	84	36905	5.779	ug/l	96
25) 2-Butanone	7.500	43	50273	17.030	ug/l	100
37) Ethyl Acetate	7.571	43	14098	2.110	ug/l #	92
43) Isopropyl Acetate	8.700	43	390423	35.499	ug/l #	78
51) 4-Methyl-2-Pentanone	10.453	43	162320	24.813	ug/l	89

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

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