

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072723\
 Data File : VN078669.D
 Acq On : 27 Jul 2023 18:58
 Operator : JC\MD
 Sample : 03768-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FL-1

Quant Time: Jul 28 01:11:10 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	404431	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	733328	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	704517	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	259577	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	261233	48.182	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	96.360%
35) Dibromofluoromethane	8.177	113	256729	51.572	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.140%
50) Toluene-d8	10.577	98	904349	48.089	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.180%
62) 4-Bromofluorobenzene	12.853	95	302621	52.243	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	104.480%
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	106453	63.031	ug/l	99
20) Methylene Chloride	5.289	84	20395	3.062	ug/l	96
25) 2-Butanone	7.494	43	62256	23.417	ug/l #	82
37) Ethyl Acetate	7.571	43	22596	3.771	ug/l #	90
43) Isopropyl Acetate	8.694	43	548997	55.651	ug/l #	71
51) 4-Methyl-2-Pentanone	10.453	43	264578	45.090	ug/l	89

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

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