

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072123\
 Data File : VN078621.D
 Acq On : 21 Jul 2023 18:13
 Operator : JC\MD
 Sample : 03680-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-3

Quant Time: Jul 21 23:28:15 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	404089	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	756491	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	737525	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	258137	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	278572	51.423	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 102.840%		
35) Dibromofluoromethane	8.177	113	280141	54.552	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 109.100%		
50) Toluene-d8	10.571	98	929768	47.927	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 95.860%		
62) 4-Bromofluorobenzene	12.853	95	316705	53.001	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 106.000%		
Target Compounds						
					Qvalue	
16) Acetone	4.441	43	68585	40.644	ug/l	94
20) Methylene Chloride	5.283	84	19903	2.962	ug/l #	92
25) 2-Butanone	7.506	43	44471	16.742	ug/l	88
37) Ethyl Acetate	7.571	43	15393	2.490	ug/l #	93
43) Isopropyl Acetate	8.694	43	299676	29.448	ug/l #	84
51) 4-Methyl-2-Pentanone	10.453	43	149405	24.682	ug/l	90

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

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