

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN073123\
 Data File : VN078707.D
 Acq On : 31 Jul 2023 15:59
 Operator : JC\MD
 Sample : 03813-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PLAUDERVILLE-2

Quant Time: Aug 01 01:19:18 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	678448	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	1246656	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	1135431	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	389156	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	450387	49.519	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.040%
35) Dibromofluoromethane	8.177	113	421488	49.805	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.620%
50) Toluene-d8	10.576	98	1557517	48.719	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	97.440%
62) 4-Bromofluorobenzene	12.853	95	512715	52.067	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	104.140%
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	171039	60.370	ug/l	94
20) Methylene Chloride	5.289	84	43802	4.272	ug/l	95
25) 2-Butanone	7.500	43	99974	22.417	ug/l	89
37) Ethyl Acetate	7.577	43	35533	3.488	ug/l #	94
43) Isopropyl Acetate	8.700	43	681919	40.662	ug/l #	79
51) 4-Methyl-2-Pentanone	10.453	43	431953	43.303	ug/l	94

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

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