

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN073123\
 Data File : VN078708.D
 Acq On : 31 Jul 2023 16:24
 Operator : JC\MD
 Sample : 03813-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PLAUDERVILLE-3

Quant Time: Aug 01 01:19:41 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	682994	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	1260113	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	1172915	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	408053	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	455088	49.702	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.400%
35) Dibromofluoromethane	8.171	113	431989	50.501	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.000%
50) Toluene-d8	10.571	98	1583887	49.015	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.020%
62) 4-Bromofluorobenzene	12.853	95	529506	53.198	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	106.400%
Target Compounds						
					Qvalue	
16) Acetone	4.453	43	172144	60.355	ug/l	94
20) Methylene Chloride	5.295	84	43367	4.180	ug/l #	89
25) 2-Butanone	7.500	43	95901	21.360	ug/l #	85
37) Ethyl Acetate	7.577	43	27698	2.690	ug/l #	95
43) Isopropyl Acetate	8.700	43	680310	40.133	ug/l #	79
51) 4-Methyl-2-Pentanone	10.453	43	431579	42.803	ug/l	92

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

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