

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN078997.D
 Acq On : 22 Aug 2023 16:24
 Operator : JC\MD
 Sample : 04090-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-I

Quant Time: Aug 23 03:04:49 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

Compound	R.T. QIon		Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.229	168	171490	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	348354	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	342145	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	107045	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	143158	53.647	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.300%	
35) Dibromofluoromethane	8.171	113	116663	49.751	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.500%	
50) Toluene-d8	10.570	98	439878	50.520	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 101.040%	
62) 4-Bromofluorobenzene	12.853	95	135930	47.763	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.520%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	25759	23.394	ug/l #	86
18) Methyl Acetate	5.041	43	5486	1.179	ug/l #	63
20) Methylene Chloride	5.271	84	11970	5.101	ug/l #	83
43) Isopropyl Acetate	8.694	43	221752	31.789	ug/l #	70
51) 4-Methyl-2-Pentanone	10.447	43	38589	9.327	ug/l #	78

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

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