

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN079003.D
 Acq On : 22 Aug 2023 18:49
 Operator : JC\MD
 Sample : 04090-11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-N

Quant Time: Aug 23 03:07:29 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	8.230	168	174973	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	344467	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	335954	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	102790	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	142188	52.223	ug/l	0.00
Spiked Amount 50.000	Range	74 - 125	Recovery	= 104.440%		
35) Dibromofluoromethane	8.171	113	121631	52.455	ug/l	0.00
Spiked Amount 50.000	Range	75 - 124	Recovery	= 104.900%		
50) Toluene-d8	10.571	98	430088	49.953	ug/l	0.00
Spiked Amount 50.000	Range	86 - 113	Recovery	= 99.900%		
62) 4-Bromofluorobenzene	12.853	95	129493	46.014	ug/l	0.00
Spiked Amount 50.000	Range	83 - 123	Recovery	= 92.020%		
Target Compounds						
						Qvalue
16) Acetone	4.436	43	26167	23.292	ug/l #	79
20) Methylene Chloride	5.283	84	13484	5.632	ug/l #	77
25) 2-Butanone	7.494	43	9244	5.132	ug/l	99
43) Isopropyl Acetate	8.694	43	187848	27.233	ug/l #	73
51) 4-Methyl-2-Pentanone	10.453	43	38927	9.515	ug/l #	80

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

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