

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082323\
 Data File : VN079016.D
 Acq On : 23 Aug 2023 15:12
 Operator : JC\MD
 Sample : 04091-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-C

Quant Time: Aug 24 01:12:02 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	172547	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	343045	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	341239	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	104595	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	150423	56.024	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 112.040%	
35) Dibromofluoromethane	8.171	113	125153	54.198	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 108.400%	
50) Toluene-d8	10.571	98	438432	51.134	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 102.260%	
62) 4-Bromofluorobenzene	12.853	95	133763	47.729	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.460%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	27129	24.488	ug/l	94
20) Methylene Chloride	5.289	84	13580	5.752	ug/l	92
25) 2-Butanone	7.500	43	9801	5.518	ug/l	99
43) Isopropyl Acetate	8.694	43	161987	23.581	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	35478	8.708	ug/l #	82

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

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