

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082323\
 Data File : VN079020.D
 Acq On : 23 Aug 2023 16:48
 Operator : JC\MD
 Sample : 04091-20
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-F

Quant Time: Aug 24 01:19:51 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	157701	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	326348	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	312386	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	89178	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	143854	58.621	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 117.240%	
35) Dibromofluoromethane	8.171	113	114951	52.326	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.660%	
50) Toluene-d8	10.571	98	414156	50.774	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 101.540%	
62) 4-Bromofluorobenzene	12.853	95	117985	44.253	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 88.500%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	28145	27.796	ug/l #	85
20) Methylene Chloride	5.271	84	13161	6.099	ug/l #	85
25) 2-Butanone	7.494	43	9052	5.576	ug/l #	71
43) Isopropyl Acetate	8.694	43	214726	32.858	ug/l #	70
51) 4-Methyl-2-Pentanone	10.453	43	38704	9.986	ug/l #	79

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

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