

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN079004.D
 Acq On : 22 Aug 2023 19:13
 Operator : JC\MD
 Sample : 04091-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-A

Quant Time: Aug 23 03:07:56 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	176401	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	354768	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	346692	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	105660	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	146983	53.547	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.100%	
35) Dibromofluoromethane	8.171	113	121788	50.998	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.000%	
50) Toluene-d8	10.571	98	444584	50.138	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 100.280%	
62) 4-Bromofluorobenzene	12.853	95	138514	47.791	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.580%	
Target Compounds						
					Qvalue	
16) Acetone	4.441	43	26546	23.438	ug/l	92
20) Methylene Chloride	5.277	84	11981	4.964	ug/l #	80
25) 2-Butanone	7.500	43	9101	5.012	ug/l #	80
43) Isopropyl Acetate	8.694	43	214818	30.239	ug/l #	70
51) 4-Methyl-2-Pentanone	10.453	43	37979	9.014	ug/l #	78

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

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