

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN078994.D
 Acq On : 22 Aug 2023 15:11
 Operator : JC\MD
 Sample : PB154968ZHE#09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154968ZHE#09

Quant Time: Aug 23 03:03:28 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	170968	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	339764	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	331071	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	96287	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	142713	53.643	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.280%	
35) Dibromofluoromethane	8.171	113	116687	51.019	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.040%	
50) Toluene-d8	10.570	98	436572	51.408	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 102.820%	
62) 4-Bromofluorobenzene	12.853	95	129376	46.609	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 93.220%	
Target Compounds						
					Qvalue	
16) Acetone	4.441	43	30420	27.712	ug/l #	84
20) Methylene Chloride	5.288	84	9944	4.251	ug/l #	87
25) 2-Butanone	7.494	43	17403	9.888	ug/l #	81
43) Isopropyl Acetate	8.694	43	213201	31.336	ug/l #	72
51) 4-Methyl-2-Pentanone	10.453	43	66227	16.412	ug/l #	80

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

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