

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN078996.D
 Acq On : 22 Aug 2023 16:00
 Operator : JC\MD
 Sample : PB154968ZHE#11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154968ZHE#11

Quant Time: Aug 23 03:04:21 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	173544	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	341624	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	342342	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	103737	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	142091	52.617	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 105.240%	
35) Dibromofluoromethane	8.171	113	115838	50.372	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.740%	
50) Toluene-d8	10.570	98	436732	51.147	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 102.300%	
62) 4-Bromofluorobenzene	12.853	95	133822	47.948	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.900%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	28772	25.822	ug/l #	78
20) Methylene Chloride	5.294	84	11504	4.845	ug/l #	80
25) 2-Butanone	7.500	43	14247	7.975	ug/l #	68
43) Isopropyl Acetate	8.694	43	234526	34.283	ug/l #	71
51) 4-Methyl-2-Pentanone	10.453	43	52657	12.978	ug/l #	82

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

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