

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN078986.D
 Acq On : 22 Aug 2023 11:58
 Operator : JC\MD
 Sample : PB154968ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154968ZHE#01

Quant Time: Aug 23 02:59:52 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	176182	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	340748	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	337955	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	101735	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	141196	51.502	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 103.000%		
35) Dibromofluoromethane	8.171	113	118389	51.614	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 103.220%		
50) Toluene-d8	10.571	98	432183	50.744	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 101.480%		
62) 4-Bromofluorobenzene	12.853	95	136171	48.915	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 97.840%		
Target Compounds						
						Qvalue
16) Acetone	4.448	43	24372	21.545	ug/l #	82
20) Methylene Chloride	5.289	84	10457	4.338	ug/l #	79
25) 2-Butanone	7.500	43	11221	6.187	ug/l #	82
43) Isopropyl Acetate	8.694	43	157442	23.074	ug/l #	79
51) 4-Methyl-2-Pentanone	10.453	43	42254	10.441	ug/l #	78

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

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