

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN078988.D
 Acq On : 22 Aug 2023 12:46
 Operator : JC\MD
 Sample : PB154968ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154968ZHE#03

Quant Time: Aug 23 03:00:48 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	175931	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	340419	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	335797	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	100297	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	144162	52.659	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 105.320%		
35) Dibromofluoromethane	8.171	113	118553	51.735	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 103.480%		
50) Toluene-d8	10.571	98	432248	50.801	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 101.600%		
62) 4-Bromofluorobenzene	12.853	95	132168	47.523	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 95.040%		
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	24399	21.600	ug/l #	78
20) Methylene Chloride	5.283	84	11227	4.664	ug/l #	86
25) 2-Butanone	7.500	43	10966	6.055	ug/l #	62
37) Ethyl Acetate	7.583	43	5147	1.335	ug/l #	86
43) Isopropyl Acetate	8.694	43	177062	25.974	ug/l #	75
51) 4-Methyl-2-Pentanone	10.447	43	39746	9.831	ug/l #	85

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

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