

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081823\
 Data File : VN078944.D
 Acq On : 18 Aug 2023 15:58
 Operator : JC\MD
 Sample : PB154902ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154902ZHE#01

Quant Time: Aug 18 23:43:33 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	158227	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	293297	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	295249	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	84128	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	124932	50.741	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 101.480%		
35) Dibromofluoromethane	8.171	113	102902	52.120	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 104.240%		
50) Toluene-d8	10.571	98	392465	53.536	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 107.080%		
62) 4-Bromofluorobenzene	12.853	95	111232	46.421	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 92.840%		
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	29683	29.218	ug/l	91
20) Methylene Chloride	5.277	84	18665	8.621	ug/l #	75
25) 2-Butanone	7.494	43	10965	6.732	ug/l #	65
43) Isopropyl Acetate	8.694	43	173827	29.597	ug/l #	79
51) 4-Methyl-2-Pentanone	10.453	43	42858	12.304	ug/l #	79

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

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