

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078829.D
 Acq On : 10 Aug 2023 11:55
 Operator : JC\MD
 Sample : PB154726TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726TB

Quant Time: Aug 10 23:48:50 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080723W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 08 05:03:52 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.230	168	162498	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	304499	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	283317	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	74561	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	125376	51.466	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 102.940%		
35) Dibromofluoromethane	8.177	113	104768	50.949	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.900%		
50) Toluene-d8	10.571	98	397585	52.540	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 105.080%		
62) 4-Bromofluorobenzene	12.859	95	103995	45.984	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 91.960%		
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	28436	30.338	ug/l #	87
20) Methylene Chloride	5.283	84	13155	5.667	ug/l #	61
25) 2-Butanone	7.494	43	14314	9.018	ug/l #	66
37) Ethyl Acetate	7.577	43	5436	1.603	ug/l #	78
43) Isopropyl Acetate	8.694	43	120742	21.114	ug/l #	80
51) 4-Methyl-2-Pentanone	10.453	43	53573	15.294	ug/l #	79

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

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