

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078833.D
 Acq On : 10 Aug 2023 13:31
 Operator : JC\MD
 Sample : PB154726ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#04

Quant Time: Aug 10 23:50:48 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	168988	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	336628	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	316208	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	88610	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	141377	55.806	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 111.620%		
35) Dibromofluoromethane	8.177	113	113685	50.009	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 100.020%		
50) Toluene-d8	10.571	98	427575	51.111	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 102.220%		
62) 4-Bromofluorobenzene	12.853	95	119358	47.740	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 95.480%		
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	32614	33.459	ug/l #	84
20) Methylene Chloride	5.283	84	14894	6.170	ug/l #	70
25) 2-Butanone	7.500	43	13887	8.413	ug/l #	70
37) Ethyl Acetate	7.577	43	5633	1.502	ug/l #	79
43) Isopropyl Acetate	8.700	43	143148	22.642	ug/l #	76
51) 4-Methyl-2-Pentanone	10.453	43	51962	13.419	ug/l #	79

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

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