

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078832.D
 Acq On : 10 Aug 2023 13:07
 Operator : JC\MD
 Sample : PB154726ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#03

Quant Time: Aug 10 23:50:21 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	8.235	168	170665	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	340301	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	306780	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	80731	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	137192	53.621	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 107.240%		
35) Dibromofluoromethane	8.171	113	112159	48.805	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 97.620%		
50) Toluene-d8	10.571	98	428701	50.692	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 101.380%		
62) 4-Bromofluorobenzene	12.853	95	113691	44.983	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 89.960%		
Target Compounds						
					Qvalue	
16) Acetone	4.453	43	31518	32.017	ug/l #	79
20) Methylene Chloride	5.294	84	13528	5.549	ug/l #	84
25) 2-Butanone	7.500	43	14612	8.766	ug/l #	77
37) Ethyl Acetate	7.583	43	5490	1.448	ug/l #	80
43) Isopropyl Acetate	8.700	43	136952	21.429	ug/l #	77
51) 4-Methyl-2-Pentanone	10.447	43	54068	13.812	ug/l #	82

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

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