

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078835.D
 Acq On : 10 Aug 2023 14:20
 Operator : JC\MD
 Sample : PB154726ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#06

Quant Time: Aug 10 23:51:45 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	8.236	168	176372	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	342948	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	319739	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	92461	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	140783	53.245	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 106.480%		
35) Dibromofluoromethane	8.177	113	118048	50.971	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.940%		
50) Toluene-d8	10.571	98	440473	51.682	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 103.360%		
62) 4-Bromofluorobenzene	12.853	95	122340	48.031	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 96.060%		
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	63126	62.050	ug/l	92
20) Methylene Chloride	5.283	84	13277	5.270	ug/l #	91
25) 2-Butanone	7.500	43	32329	18.766	ug/l #	78
37) Ethyl Acetate	7.577	43	32265	8.446	ug/l #	94
43) Isopropyl Acetate	8.700	43	178692	27.744	ug/l #	75
51) 4-Methyl-2-Pentanone	10.453	43	110508	28.011	ug/l #	81

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

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