

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078836.D
 Acq On : 10 Aug 2023 14:44
 Operator : JC\MD
 Sample : PB154726ZHE#07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#07

Quant Time: Aug 10 23:52:06 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	163896	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	328708	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	306143	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	85405	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	131002	53.317	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 106.640%			
35) Dibromofluoromethane	8.171	113	109799	49.463	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 98.920%			
50) Toluene-d8	10.577	98	432285	52.919	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 105.840%			
62) 4-Bromofluorobenzene	12.859	95	113903	46.656	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 93.320%			
Target Compounds						
16) Acetone	4.442	43	62663	66.284	ug/l	99
20) Methylene Chloride	5.295	84	13660	5.834	ug/l #	83
25) 2-Butanone	7.494	43	32886	20.543	ug/l #	72
37) Ethyl Acetate	7.571	43	32819	8.963	ug/l #	93
43) Isopropyl Acetate	8.694	43	191329	30.993	ug/l #	73
51) 4-Methyl-2-Pentanone	10.453	43	112080	29.641	ug/l #	84

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

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