

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN093023\
 Data File : VN079385.D
 Acq On : 30 Sep 2023 14:34
 Operator : JC\MD
 Sample : PB155992ZHE#05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155992ZHE#05

Quant Time: Oct 02 01:55:35 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091123W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 12 05:05:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	127260	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	260834	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	264796	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	66809	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	113598	66.127	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 132.260%#	
35) Dibromofluoromethane	8.171	113	95168	54.851	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 109.700%	
50) Toluene-d8	10.571	98	350548	55.025	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 110.040%	
62) 4-Bromofluorobenzene	12.853	95	95507	46.998	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 94.000%	
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	21232	29.101	ug/l #	81
20) Methylene Chloride	5.271	84	13722	7.729	ug/l #	87
25) 2-Butanone	7.500	43	5487	5.183	ug/l #	85
43) Isopropyl Acetate	8.694	43	146085	33.564	ug/l #	76
51) 4-Methyl-2-Pentanone	10.447	43	14164	5.606	ug/l #	81

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

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