

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN093023\
 Data File : VN079381.D
 Acq On : 30 Sep 2023 12:58
 Operator : JC\MD
 Sample : PB155992ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155992ZHE#01

Quant Time: Oct 02 01:53:31 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	112262	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	221494	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	218038	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	50621	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	99677	65.775	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 131.540%#	
35) Dibromofluoromethane	8.177	113	81283	55.169	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 110.340%	
50) Toluene-d8	10.571	98	298597	55.195	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 110.380%	
62) 4-Bromofluorobenzene	12.853	95	76475	44.317	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 88.640%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	23025	35.775	ug/l	96
20) Methylene Chloride	5.283	84	12461	7.956	ug/l #	83
25) 2-Butanone	7.500	43	5635	6.034	ug/l #	66
37) Ethyl Acetate	7.577	43	3741	1.652	ug/l #	78
43) Isopropyl Acetate	8.694	43	128434	34.750	ug/l #	81
51) 4-Methyl-2-Pentanone	10.453	43	16211	7.556	ug/l #	71

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

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