

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN093023\
 Data File : VN079382.D
 Acq On : 30 Sep 2023 13:22
 Operator : JC\MD
 Sample : PB155992ZHE#02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155992ZHE#02

Quant Time: Oct 02 01:54:07 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	141403	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	289628	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	291967	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	75737	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	125492	65.744	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 131.480%#			
35) Dibromofluoromethane	8.171	113	106591	55.327	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 110.660%			
50) Toluene-d8	10.571	98	389621	55.078	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 110.160%			
62) 4-Bromofluorobenzene	12.853	95	108581	48.120	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 96.240%			
Target Compounds						Qvalue
16) Acetone	4.436	43	23388	28.850	ug/l #	84
20) Methylene Chloride	5.277	84	14104	7.149	ug/l	86
37) Ethyl Acetate	7.571	43	4002	1.351	ug/l #	79
43) Isopropyl Acetate	8.694	43	132803	27.479	ug/l #	80
51) 4-Methyl-2-Pentanone	10.447	43	16564	5.904	ug/l #	70

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

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