

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN093023\
 Data File : VN079384.D
 Acq On : 30 Sep 2023 14:10
 Operator : JC\MD
 Sample : PB155992ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155992ZHE#04

Quant Time: Oct 02 01:55:06 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	8.229	168	132790	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	270776	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	273130	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	72473	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	117614	65.613	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 131.220%#	
35) Dibromofluoromethane	8.177	113	96288	53.459	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 106.920%	
50) Toluene-d8	10.571	98	364683	55.142	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 110.280%	
62) 4-Bromofluorobenzene	12.853	95	97961	46.436	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 92.880%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	21715	28.524	ug/l #	77
20) Methylene Chloride	5.283	84	12866	6.945	ug/l #	83
43) Isopropyl Acetate	8.694	43	139894	30.962	ug/l #	77
51) 4-Methyl-2-Pentanone	10.453	43	14592	5.563	ug/l #	70

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

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