

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091523\
 Data File : VN079279.D
 Acq On : 15 Sep 2023 18:07
 Operator : JC\MD
 Sample : PB155573ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155573ZHE#03

Quant Time: Sep 16 01:56:51 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	133095	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	272355	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	271674	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	75991	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	116275	64.717	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 129.440%#	
35) Dibromofluoromethane	8.171	113	95800	52.880	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 105.760%	
50) Toluene-d8	10.571	98	359648	54.065	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 108.140%	
62) 4-Bromofluorobenzene	12.853	95	105349	49.649	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 99.300%	
Target Compounds						
					Qvalue	
16) Acetone	4.453	43	18447	24.175	ug/l	89
20) Methylene Chloride	5.283	84	11965	6.444	ug/l #	85
25) 2-Butanone	7.500	43	11232	10.145	ug/l #	84
43) Isopropyl Acetate	8.694	43	172721	38.005	ug/l #	75
51) 4-Methyl-2-Pentanone	10.447	43	30353	11.505	ug/l #	79

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

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