

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091523\
 Data File : VN079273.D
 Acq On : 15 Sep 2023 15:42
 Operator : JC\MD
 Sample : PB155573ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155573ZHE#04

Quant Time: Sep 16 01:43:59 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	151650	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	299574	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	300489	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	91420	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	124249	60.694	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 121.380%			
35) Dibromofluoromethane	8.171	113	101846	51.109	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 102.220%			
50) Toluene-d8	10.571	98	398428	54.453	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 108.900%			
62) 4-Bromofluorobenzene	12.853	95	120236	51.516	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 103.040%			
Target Compounds						
16) Acetone	4.436	43	19659	22.611	ug/l #	78
20) Methylene Chloride	5.283	84	12011	5.677	ug/l #	74
25) 2-Butanone	7.494	43	12512	9.919	ug/l #	73
43) Isopropyl Acetate	8.694	43	162887	32.585	ug/l #	79
51) 4-Methyl-2-Pentanone	10.453	43	34395	11.853	ug/l #	77

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

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