

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091523\
 Data File : VN079270.D
 Acq On : 15 Sep 2023 14:30
 Operator : JC\MD
 Sample : PB155573ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155573ZHE#01

Quant Time: Sep 16 01:41:44 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	151108	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	293119	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	286326	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	92278	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	119562	58.614	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 117.220%		
35) Dibromofluoromethane	8.171	113	101915	52.270	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 104.540%		
50) Toluene-d8	10.570	98	381094	53.231	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 106.460%		
62) 4-Bromofluorobenzene	12.853	95	113953	49.899	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 99.800%		
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	92916	107.254	ug/l	96
20) Methylene Chloride	5.277	84	17521	8.311	ug/l #	80
25) 2-Butanone	7.488	43	21428	17.047	ug/l	92
43) Isopropyl Acetate	8.694	43	152105	31.098	ug/l #	81
51) 4-Methyl-2-Pentanone	10.453	43	31104	10.955	ug/l #	80

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

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