

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110623\
 Data File : VN079917.D
 Acq On : 06 Nov 2023 12:58
 Operator : JC\MD
 Sample : PB156885TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB156885TB

Quant Time: Nov 07 04:45:52 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110123W.M
 Quant Title : SW846 8260
 QLast Update : Thu Nov 02 05:49:08 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.229	168	262684	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	525021	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	564058	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	212290	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	219981	54.967	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 109.940%			
35) Dibromofluoromethane	8.171	113	181607	48.873	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 97.740%			
50) Toluene-d8	10.571	98	675420	49.375	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 98.740%			
62) 4-Bromofluorobenzene	12.853	95	234357	51.657	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 103.320%			
Target Compounds						Qvalue
16) Acetone	4.430	43	102212	72.709	ug/l	90
20) Methylene Chloride	5.271	84	21805	6.147	ug/l #	77
25) 2-Butanone	7.488	43	10582	5.141	ug/l #	64
37) Ethyl Acetate	7.571	43	6945	1.427	ug/l #	83
43) Isopropyl Acetate	8.694	43	227096	26.613	ug/l #	80
51) 4-Methyl-2-Pentanone	10.447	43	24916	5.186	ug/l #	76

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

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