

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110623\
 Data File : VN079933.D
 Acq On : 06 Nov 2023 19:21
 Operator : JC\MD
 Sample : PB156885ZHE#16
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB156885ZHE#16

Quant Time: Nov 07 04:53:22 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	8.230	168	232727	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	512584	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	553579	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	198251	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	225822	63.689	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 127.380%#		
35) Dibromofluoromethane	8.171	113	188748	52.027	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 104.060%		
50) Toluene-d8	10.571	98	707642	52.985	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 105.980%		
62) 4-Bromofluorobenzene	12.853	95	234222	52.880	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 105.760%		
Target Compounds						
						Qvalue
16) Acetone	4.436	43	60297	48.414	ug/l #	88
18) Methyl Acetate	5.042	43	10864	1.823	ug/l #	82
20) Methylene Chloride	5.277	84	10565	3.362	ug/l #	83
37) Ethyl Acetate	7.577	43	5929	1.248	ug/l #	79
43) Isopropyl Acetate	8.694	43	351491	42.190	ug/l #	74
51) 4-Methyl-2-Pentanone	10.447	43	23536	5.018	ug/l #	65

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

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