

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110623\
 Data File : VN079923.D
 Acq On : 06 Nov 2023 15:21
 Operator : JC\MD
 Sample : PB156885ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB156885ZHE#06

Quant Time: Nov 07 04:48:46 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	229907	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	501566	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	534508	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	188456	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	220562	62.969	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 125.940%#	
35) Dibromofluoromethane	8.171	113	181180	51.038	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.080%	
50) Toluene-d8	10.570	98	681302	52.134	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 104.260%	
62) 4-Bromofluorobenzene	12.853	95	222372	51.307	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 102.620%	
Target Compounds						
						Qvalue
16) Acetone	4.430	43	65470	53.212	ug/l	90
20) Methylene Chloride	5.283	84	12207	3.932	ug/l	93
37) Ethyl Acetate	7.565	43	5254	1.130	ug/l #	73
43) Isopropyl Acetate	8.694	43	350485	42.993	ug/l #	74
51) 4-Methyl-2-Pentanone	10.453	43	23865	5.200	ug/l #	65

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

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