

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110623\
 Data File : VN079928.D
 Acq On : 06 Nov 2023 17:21
 Operator : JC\MD
 Sample : PB156885ZHE#11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB156885ZHE#11

Quant Time: Nov 07 04:51:07 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	221671	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	493426	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	529417	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	199196	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	214518	63.519	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 127.040%#		
35) Dibromofluoromethane	8.171	113	177298	50.769	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.540%		
50) Toluene-d8	10.570	98	675073	52.509	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 105.020%		
62) 4-Bromofluorobenzene	12.853	95	231775	54.359	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 108.720%		
Target Compounds						
					Qvalue	
16) Acetone	4.441	43	62430	52.626	ug/l	88
18) Methyl Acetate	5.041	43	9659	1.702	ug/l #	82
20) Methylene Chloride	5.283	84	12917	4.315	ug/l #	78
37) Ethyl Acetate	7.577	43	6410	1.402	ug/l #	83
43) Isopropyl Acetate	8.694	43	350575	43.713	ug/l #	73
51) 4-Methyl-2-Pentanone	10.447	43	23499	5.205	ug/l #	63

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

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