

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110623\
 Data File : VN079921.D
 Acq On : 06 Nov 2023 14:33
 Operator : JC\MD
 Sample : PB156885ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB156885ZHE#04

Quant Time: Nov 07 04:47:48 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	248576	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	534833	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	578768	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	218244	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	229268	60.539	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 121.080%		
35) Dibromofluoromethane	8.171	113	190573	50.345	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 100.700%		
50) Toluene-d8	10.571	98	725709	52.078	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 104.160%		
62) 4-Bromofluorobenzene	12.853	95	250989	54.308	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 108.620%		
Target Compounds						
						Qvalue
16) Acetone	4.430	43	105097	79.004	ug/l	94
18) Methyl Acetate	5.047	43	9875	1.552	ug/l #	68
20) Methylene Chloride	5.283	84	21996	6.553	ug/l #	82
25) 2-Butanone	7.500	43	11590	5.950	ug/l #	74
37) Ethyl Acetate	7.583	43	6845	1.381	ug/l #	86
43) Isopropyl Acetate	8.694	43	286597	32.969	ug/l #	76
51) 4-Methyl-2-Pentanone	10.447	43	24786	5.065	ug/l #	71

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

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