

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020724\
 Data File : VN080943.D
 Acq On : 07 Feb 2024 18:03
 Operator : JC\MD
 Sample : PB158850ZHE#07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB158850ZHE#07

Quant Time: Feb 08 00:29:36 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N020624W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 07 06:22:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	267443	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	9.106	114	588920	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	738404	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	313311	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	190206	56.377	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 112.760%	
35) Dibromofluoromethane	8.171	113	175299	55.472	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 110.940%	
50) Toluene-d8	10.571	98	697984	61.063	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 122.120%#	
62) 4-Bromofluorobenzene	12.853	95	299692	72.173	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 144.340%#	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	45876	41.331	ug/l	91
20) Methylene Chloride	5.277	84	10043	2.986	ug/l	# 79
37) Ethyl Acetate	7.559	43	7568	2.127	ug/l	# 88
43) Isopropyl Acetate	8.694	43	304528	46.554	ug/l	# 90
51) 4-Methyl-2-Pentanone	10.447	43	20812	5.279	ug/l	# 79

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

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