

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082433.D
 Acq On : 03 Jun 2024 14:47
 Operator : JC\MD
 Sample : PB161254ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161254ZHE#01

Quant Time: Jun 04 00:29:14 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 16 05:21:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	165592	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	308436	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	307815	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	133306	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	143640	54.935	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 109.860%		
35) Dibromofluoromethane	8.165	113	102274	53.903	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 107.800%		
50) Toluene-d8	10.571	98	405195	55.702	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 111.400%		
62) 4-Bromofluorobenzene	12.853	95	141153	54.365	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	= 108.740%		
Target Compounds						
						Qvalue
16) Acetone	4.436	43	84577	78.851	ug/l	90
20) Methylene Chloride	5.277	84	12619	5.831	ug/l #	83
25) 2-Butanone	7.494	43	14400	8.855	ug/l #	84
43) Isopropyl Acetate	8.694	43	336475	53.322	ug/l #	70
51) 4-Methyl-2-Pentanone	10.447	43	20374	5.683	ug/l #	72

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

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