

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082442.D
 Acq On : 03 Jun 2024 18:27
 Operator : JC\MD
 Sample : PB161254ZHE#09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161254ZHE#09

Quant Time: Jun 04 00:33:47 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	8.224	168	155027	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	284495	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	289004	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	124403	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	135778	55.467	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 110.940%		
35) Dibromofluoromethane	8.171	113	96100	54.912	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 109.820%		
50) Toluene-d8	10.571	98	376617	56.130	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 112.260%		
62) 4-Bromofluorobenzene	12.853	95	132456	55.309	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	= 110.620%		
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	86031	85.673	ug/l	93
20) Methylene Chloride	5.271	84	12917	6.375	ug/l #	86
25) 2-Butanone	7.489	43	13992	9.190	ug/l #	78
43) Isopropyl Acetate	8.694	43	306556	52.669	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	20657	6.246	ug/l #	72

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

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