

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082437.D
 Acq On : 03 Jun 2024 16:25
 Operator : JC\MD
 Sample : PB161254ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161254ZHE#04

Quant Time: Jun 04 00:31:28 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	157845	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	295963	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	292650	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	120678	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	136505	54.768	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 109.540%		
35) Dibromofluoromethane	8.171	113	98845	54.292	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 108.580%		
50) Toluene-d8	10.571	98	387168	55.467	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 110.940%		
62) 4-Bromofluorobenzene	12.853	95	134258	53.889	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	= 107.780%		
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	84816	82.955	ug/l	96
20) Methylene Chloride	5.283	84	13219	6.408	ug/l #	87
25) 2-Butanone	7.494	43	13858	8.940	ug/l #	80
43) Isopropyl Acetate	8.694	43	299988	49.543	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	20247	5.885	ug/l #	71

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

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