

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082435.D
 Acq On : 03 Jun 2024 15:36
 Operator : JC\MD
 Sample : PB161254ZHE#02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161254ZHE#02

Quant Time: Jun 04 00:30:30 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	161147	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	296739	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	295612	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	125037	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	137962	54.219	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 108.440%	
35) Dibromofluoromethane	8.165	113	100387	54.994	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 109.980%	
50) Toluene-d8	10.571	98	388429	55.502	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 111.000%	
62) 4-Bromofluorobenzene	12.853	95	135967	54.432	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	= 108.860%	
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	84826	81.265	ug/l	90
20) Methylene Chloride	5.277	84	12921	6.135	ug/l #	78
25) 2-Butanone	7.494	43	13685	8.647	ug/l #	86
43) Isopropyl Acetate	8.688	43	296006	48.758	ug/l #	73
51) 4-Methyl-2-Pentanone	10.447	43	21089	6.114	ug/l #	71

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

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