

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN053124\
 Data File : VN082407.D
 Acq On : 31 May 2024 14:46
 Operator : JC\MD
 Sample : PB161190ZHE#05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161190ZHE#05

Quant Time: May 31 22:58:21 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	157279	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	287806	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	285518	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	116040	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	133338	53.690	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 107.380%		
35) Dibromofluoromethane	8.165	113	95242	53.795	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 107.600%		
50) Toluene-d8	10.571	98	377166	55.566	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 111.140%		
62) 4-Bromofluorobenzene	12.847	95	128364	52.984	ug/l	0.00
Spiked Amount	50.000	Range 64 - 133	Recovery	= 105.960%		
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	73972	72.609	ug/l	91
20) Methylene Chloride	5.283	84	12323	5.995	ug/l #	84
25) 2-Butanone	7.488	43	11385	7.371	ug/l	99
43) Isopropyl Acetate	8.688	43	329788	56.008	ug/l #	69
51) 4-Methyl-2-Pentanone	10.447	43	18736	5.600	ug/l #	70

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

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