

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN053124\
 Data File : VN082409.D
 Acq On : 31 May 2024 15:34
 Operator : JC\MD
 Sample : PB161190ZHE#07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161190ZHE#07

Quant Time: May 31 22:59:12 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	156579	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	290560	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	291308	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	121636	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	135051	54.623	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 109.240%	
35) Dibromofluoromethane	8.171	113	96566	54.026	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 108.060%	
50) Toluene-d8	10.571	98	386680	56.427	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 112.860%	
62) 4-Bromofluorobenzene	12.847	95	133155	54.440	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	= 108.880%	
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	75354	74.297	ug/l	92
20) Methylene Chloride	5.277	84	12046	5.887	ug/l #	82
25) 2-Butanone	7.494	43	12467	8.107	ug/l #	82
43) Isopropyl Acetate	8.688	43	290793	48.918	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	18461	5.466	ug/l #	72

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

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