

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN053124\
 Data File : VN082405.D
 Acq On : 31 May 2024 13:58
 Operator : JC\MD
 Sample : PB161190ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161190ZHE#03

Quant Time: May 31 22:57:33 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	158107	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	293346	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	290009	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	125945	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	136442	54.652	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 109.300%	
35) Dibromofluoromethane	8.165	113	97481	54.020	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 108.040%	
50) Toluene-d8	10.571	98	389358	56.278	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 112.560%	
62) 4-Bromofluorobenzene	12.847	95	133220	53.949	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	= 107.900%	
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	74828	73.065	ug/l	90
20) Methylene Chloride	5.277	84	13597	6.580	ug/l #	77
25) 2-Butanone	7.488	43	11967	7.707	ug/l	89
43) Isopropyl Acetate	8.688	43	290921	48.474	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	18847	5.527	ug/l #	72

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

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