

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
 Data File : VN082637.D
 Acq On : 19 Jun 2024 13:33
 Operator : JC\MD
 Sample : PB161542ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161542ZHE#03

Quant Time: Jun 20 03:00:11 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061724W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jun 18 04:52:44 2024
 Response via : Initial Calibration

Compound	R.T. QIon		Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.230	168	129565	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	247496	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	259494	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	101460	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	118489	57.028	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 114.060%	
35) Dibromofluoromethane	8.171	113	89555	53.339	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 106.680%	
50) Toluene-d8	10.571	98	334811	52.732	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 105.460%	
62) 4-Bromofluorobenzene	12.853	95	109220	49.894	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	= 99.780%	
Target Compounds						
						Qvalue
16) Acetone	4.442	43	59182	62.631	ug/l	98
20) Methylene Chloride	5.283	84	11740	7.067	ug/l	95
25) 2-Butanone	7.489	43	17122	12.748	ug/l	98
43) Isopropyl Acetate	8.694	43	251739	42.132	ug/l #	82
51) 4-Methyl-2-Pentanone	10.453	43	15996	5.220	ug/l #	79

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

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