

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061924\
 Data File : VN082640.D
 Acq On : 19 Jun 2024 14:47
 Operator : JC\MD
 Sample : PB161542ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161542ZHE#06

Quant Time: Jun 20 03:01:38 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	132229	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	254151	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	267131	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	105391	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	121770	57.426	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	114.860%
35) Dibromofluoromethane	8.171	113	89958	52.176	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	104.360%
50) Toluene-d8	10.571	98	343484	52.681	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	105.360%
62) 4-Bromofluorobenzene	12.853	95	110514	49.163	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	=	98.320%
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	58076	60.222	ug/l	100
20) Methylene Chloride	5.283	84	11435	6.745	ug/l	94
25) 2-Butanone	7.488	43	19131	13.957	ug/l	95
43) Isopropyl Acetate	8.694	43	271952	44.323	ug/l #	80
51) 4-Methyl-2-Pentanone	10.447	43	16287	5.175	ug/l #	79

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

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