

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060524\
 Data File : VN082466.D
 Acq On : 05 Jun 2024 13:24
 Operator : JC\MD
 Sample : PB161297ZHE#02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161297ZHE#02

Manual Integrations
 APPROVED

Reviewed By : John Carlone 06/06/2024
 Supervised By : Mahesh Dadoda 06/06/2024

Quant Time: Jun 06 00:55:28 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	154687	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	288968	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	291196	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	119109	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	135999	55.679	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 111.360%			
35) Dibromofluoromethane	8.171	113	98991	55.688	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 111.380%			
50) Toluene-d8	10.571	98	386512	56.713	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 113.420%#			
62) 4-Bromofluorobenzene	12.853	95	129234	53.128	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 106.260%			
Target Compounds						
16) Acetone	4.436	43	97284	97.092	ug/l	93
20) Methylene Chloride	5.283	84	4224	2.089	ug/l #	80
25) 2-Butanone	7.494	43	24453	16.096	ug/l #	84
43) Isopropyl Acetate	8.694	43	258768	43.770	ug/l #	73
51) 4-Methyl-2-Pentanone	10.447	43	23806m	7.087	ug/l	

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

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