

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082439.D
 Acq On : 03 Jun 2024 17:13
 Operator : JC\MD
 Sample : PB161254ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161254ZHE#06

Manual Integrations
 APPROVED

Reviewed By : John Carlone 06/04/2024
 Supervised By : Mahesh Dadoda 06/04/2024

Quant Time: Jun 04 00:32:23 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	158854	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	297626	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	295736	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	125997	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	138967	55.402	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 110.800%			
35) Dibromofluoromethane	8.171	113	98873	54.004	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 108.000%			
50) Toluene-d8	10.571	98	391631	55.793	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 111.580%			
62) 4-Bromofluorobenzene	12.853	95	135656	54.146	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 108.300%			
Target Compounds						
16) Acetone	4.436	43	99637	96.832	ug/l	93
20) Methylene Chloride	5.271	84	3229	1.555	ug/l #	76
25) 2-Butanone	7.494	43	25633	16.431	ug/l #	81
43) Isopropyl Acetate	8.694	43	302134	49.619	ug/l #	71
51) 4-Methyl-2-Pentanone	10.447	43	25159m	7.272	ug/l	

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

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