

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060524\
 Data File : VN082467.D
 Acq On : 05 Jun 2024 13:48
 Operator : JC\MD
 Sample : PB161297ZHE#03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161297ZHE#03

Manual Integrations
 APPROVED

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

Quant Time: Jun 06 00:55:54 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	8.229	168	158611	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	295352	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	299316	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	122379	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	138976	55.490	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 110.980%			
35) Dibromofluoromethane	8.171	113	100722	55.437	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 110.880%			
50) Toluene-d8	10.571	98	394048	56.569	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 113.140%#			
62) 4-Bromofluorobenzene	12.853	95	133255	53.597	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 107.200%			
Target Compounds					Qvalue	
16) Acetone	4.436	43	98731	96.099	ug/l	90
20) Methylene Chloride	5.277	84	4254m	2.052	ug/l	
25) 2-Butanone	7.488	43	24379	15.651	ug/l	90
43) Isopropyl Acetate	8.688	43	281120	46.523	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	19508m	5.682	ug/l	

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

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