

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082440.D
 Acq On : 03 Jun 2024 17:38
 Operator : JC\MD
 Sample : PB161254ZHE#07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161254ZHE#07

Manual Integrations
 APPROVED

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

Quant Time: Jun 04 00:32:50 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	150745	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	279041	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	274349	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	109699	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	131197	55.118	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 110.240%			
35) Dibromofluoromethane	8.171	113	92865	54.100	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 108.200%			
50) Toluene-d8	10.571	98	365217	55.495	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 111.000%			
62) 4-Bromofluorobenzene	12.853	95	120436	51.273	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 102.540%			
Target Compounds						
16) Acetone	4.430	43	100727	103.157	ug/l	90
20) Methylene Chloride	5.289	84	3413	1.732	ug/l #	85
25) 2-Butanone	7.489	43	25308	17.095	ug/l	90
43) Isopropyl Acetate	8.694	43	291767	51.108	ug/l #	72
51) 4-Methyl-2-Pentanone	10.453	43	25609m	7.895	ug/l	

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

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