

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060324\
 Data File : VN082441.D
 Acq On : 03 Jun 2024 18:02
 Operator : JC\MD
 Sample : PB161254ZHE#08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB161254ZHE#08

Manual Integrations
 APPROVED

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

Quant Time: Jun 04 00:33:19 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	156721	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	291887	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	290456	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	121221	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	138398	55.926	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 111.860%			
35) Dibromofluoromethane	8.171	113	100104	55.751	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 111.500%			
50) Toluene-d8	10.571	98	389695	56.609	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 113.220%#			
62) 4-Bromofluorobenzene	12.853	95	131371	53.467	ug/l	0.00
Spiked Amount 50.000	Range 64 - 133		Recovery = 106.940%			
Target Compounds					Qvalue	
16) Acetone	4.436	43	102142	100.618	ug/l	99
20) Methylene Chloride	5.283	84	3230	1.577	ug/l #	91
25) 2-Butanone	7.488	43	27170	17.653	ug/l #	79
37) Ethyl Acetate	7.571	43	5233	1.553	ug/l #	79
43) Isopropyl Acetate	8.694	43	342816	57.407	ug/l #	69
51) 4-Methyl-2-Pentanone	10.453	43	24322m	7.168	ug/l	

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

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