

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030124\
 Data File : VN081284.D
 Acq On : 01 Mar 2024 14:44
 Operator : JC\MD
 Sample : PB153949ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB153949ZHE#06

Manual Integrations
 APPROVED

Reviewed By : John Carlone 03/04/2024
 Supervised By : Mahesh Dadoda 03/04/2024

Quant Time: Mar 01 22:18:12 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021624W.M
 Quant Title : SW846 8260
 QLast Update : Fri Feb 16 23:45:20 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.224	168	306981	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	583781	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	529036	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	207825	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	230961	61.381	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 122.760%			
35) Dibromofluoromethane	8.165	113	179055	54.241	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 108.480%			
50) Toluene-d8	10.565	98	708392	60.419	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 120.840%#			
62) 4-Bromofluorobenzene	12.847	95	223736	50.785	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 101.560%			
Target Compounds						
16) Acetone	4.430	43	71549	48.326	ug/l	97
18) Methyl Acetate	5.030	43	4798	1.328	ug/l #	71
20) Methylene Chloride	5.283	84	35179	8.412	ug/l #	85
37) Ethyl Acetate	7.559	43	7124m	1.396	ug/l	
43) Isopropyl Acetate	8.688	43	504902	56.560	ug/l #	74

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

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