

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091024\
 Data File : VN083770.D
 Acq On : 11 Sep 2024 04:27
 Operator : JC\MD
 Sample : PB163204TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB163204TB

Manual Integrations
 APPROVED

Reviewed By :John Carlone 09/11/2024
 Supervised By :Semsettin Yesilyurt 09/11/2024

Quant Time: Sep 11 06:22:14 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091024W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 11 03:05:46 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	110570	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	232747	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	251077	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	116130	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	87127	52.258	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 104.520%			
35) Dibromofluoromethane	8.165	113	68780	46.161	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 92.320%			
50) Toluene-d8	10.565	98	259119	50.304	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 100.600%			
62) 4-Bromofluorobenzene	12.847	95	112637	53.001	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 106.000%			
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	19999	29.923	ug/l	99
20) Methylene Chloride	5.259	84	5535	4.047	ug/l	91
37) Ethyl Acetate	7.565	43	3125	1.535	ug/l #	83
43) Isopropyl Acetate	8.688	43	182532m	49.481	ug/l	

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

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