

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN051925\
 Data File : VN086687.D
 Acq On : 19 May 2025 14:15
 Operator : JC\MD
 Sample : PB168034ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB168034ZHE#04

Manual Integrations
 APPROVED

Reviewed By : John Carlone 05/21/2025
 Supervised By : Mahesh Dadoda 05/21/2025

Quant Time: May 20 03:18:51 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051625W.M
 Quant Title : SW846 8260
 QLast Update : Sat May 17 00:46:13 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	203868	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	389357	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	383243	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.788	152	157036	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.577	65	140972	50.042	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 100.080%			
35) Dibromofluoromethane	8.165	113	116465	50.227	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.460%			
50) Toluene-d8	10.565	98	509433	53.015	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 106.040%			
62) 4-Bromofluorobenzene	12.847	95	181608	53.525	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 107.060%			
Target Compounds						
16) Acetone	4.424	43	24673	23.414	ug/l	98
18) Methyl Acetate	5.036	43	3889	1.319	ug/l #	73
20) Methylene Chloride	5.277	84	16142	5.805	ug/l	95
25) 2-Butanone	7.489	43	9765	5.997	ug/l	95
43) Isopropyl Acetate	8.683	43	40158m	5.798	ug/l	

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

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