

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN092625\
 Data File : VN087966.D
 Acq On : 26 Sep 2025 15:06
 Operator : JC\MD
 Sample : PB169841TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB169841TB

Manual Integrations
 APPROVED

Reviewed By :Mahesh Dadoda 09/29/2025
 Supervised By :Semsettin Yesilyurt 09/29/2025

Quant Time: Sep 27 01:26:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N092325W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 24 05:05:30 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	183387	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	416895	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	441917	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	196760	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	173710	56.285	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 112.560%			
35) Dibromofluoromethane	8.147	113	151413	50.023	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.040%			
50) Toluene-d8	10.547	98	517622	48.629	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 97.260%			
62) 4-Bromofluorobenzene	12.829	95	176497	46.388	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 92.780%			
Target Compounds					Qvalue	
16) Acetone	4.424	43	76149	59.968	ug/l	97
18) Methyl Acetate	5.018	43	10469	3.294	ug/l	93
20) Methylene Chloride	5.253	84	11039m	4.295	ug/l	
43) Isopropyl Acetate	8.683	43	22618	3.006	ug/l #	90

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

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