

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072325\
 Data File : VX047098.D
 Acq On : 23 Jul 2025 17:36
 Operator : JC/MD
 Sample : PB168966TB
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB168966TB

Manual Integrations
 APPROVED

Reviewed By : John Carlone 07/24/2025
 Supervised By : Mahesh Dadoda 07/24/2025

Quant Time: Jul 24 05:55:21 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072125W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jul 22 03:59:51 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	479643	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	866424	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	802189	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	393927	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	314539	52.644	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 105.280%			
35) Dibromofluoromethane	5.403	113	257775	48.186	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 96.380%			
50) Toluene-d8	8.647	98	939665	54.239	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 108.480%			
62) 4-Bromofluorobenzene	11.079	95	374821	50.204	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 100.400%			
Target Compounds						
16) Acetone	2.386	43	57559	24.048	ug/l	98
18) Methyl Acetate	2.715	43	31199	4.905	ug/l	100
20) Methylene Chloride	2.806	84	24625	3.740	ug/l	99
37) Ethyl Acetate	4.745	43	18609m	2.215	ug/l	
43) Isopropyl Acetate	6.354	43	14641	1.118	ug/l	96

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

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