

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092325\
 Data File : VX047760.D
 Acq On : 24 Sep 2025 08:22
 Operator : JC/MD
 Sample : Q3141-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-VPB-184-GW-100-102

Quant Time: Sep 24 08:43:27 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091625W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 17 06:39:58 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.537	168	165497	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	300349	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	276860	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	136496	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	149951	56.923	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 113.840%			
35) Dibromofluoromethane	5.373	113	113994	56.592	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 113.180%			
50) Toluene-d8	8.634	98	383307	54.995	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 109.980%			
62) 4-Bromofluorobenzene	11.067	95	148783	56.246	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 112.500%			
Target Compounds					Qvalue	
16) Acetone	2.373	43	12472	10.889	ug/l	99
30) Chloroform	5.086	83	2823	0.668	ug/l	95
44) Trichloroethene	7.123	130	2022	0.936	ug/l	88
47) Bromodichloromethane	7.805	83	1217	0.354	ug/l #	94
52) Toluene	8.714	92	1845	0.335	ug/l	98
60) Dibromochloromethane	9.512	129	1095	0.447	ug/l	96
64) Tetrachloroethene	9.262	164	474	0.263	ug/l #	87

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

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