

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052825\
 Data File : VX046383.D
 Acq On : 28 May 2025 14:16
 Operator : JC/MD
 Sample : IBLK
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 IBLK

Quant Time: May 29 03:36:55 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X050525W.M
 Quant Title : SW846 8260
 QLast Update : Tue May 06 07:12:22 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	58479	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	116719	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	111708	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	47222	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	60067	55.095	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 110.200%			
35) Dibromofluoromethane	5.379	113	43226	51.429	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 102.860%			
50) Toluene-d8	8.647	98	149950	51.545	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 103.100%			
62) 4-Bromofluorobenzene	11.079	95	56387	50.531	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 101.060%			
Target Compounds			Qvalue			
3) Chloromethane	1.301	50	19	0.022	ug/l #	43
6) Chloroethane	1.618	64	27	0.063	ug/l #	44
17) Carbon Disulfide	2.502	76	46	0.028	ug/l #	76
23) Vinyl Acetate	3.721	43	21	0.009	ug/l #	74
32) 1,1,1-Trichloroethane	5.507	97	21	0.016	ug/l #	22
51) 4-Methyl-2-Pentanone	8.586	43	186	0.132	ug/l #	39
59) 2-Hexanone	9.439	43	21	0.020	ug/l #	28
67) Ethyl Benzene	10.055	91	257	0.060	ug/l #	45
73) Isopropylbenzene	11.079	105	85	0.023	ug/l #	49
74) N-amyl acetate	11.073	43	83	0.046	ug/l #	1
75) 1,1,2,2-Tetrachloroethane	11.079	83	23	0.018	ug/l #	1
78) n-propylbenzene	11.085	91	294	0.069	ug/l #	54

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

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