

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060524\
 Data File : VU059150.D
 Acq On : 05 Jun 2024 14:00
 Operator : MD/SY
 Sample : VIBLK
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VIBLK

Quant Time: Jun 06 03:31:48 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U060524DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Thu Jun 06 03:28:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Fluorobenzene	6.110	96	41353	1.000	ug/l	# 0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.630	95	12322	0.876	ug/l	0.00
Spiked Amount 1.000			Recovery	=	88.000%	
68) 1,2-Dichlorobenzene-d4	12.193	152	12743	0.865	ug/l	0.00
Spiked Amount 1.000			Recovery	=	87.000%	
Target Compounds						
						Qvalue
15) Methylene Chloride	3.039	84	8558	0.627	ug/l	95
75) 1,3,5-Trimethylbenzene	11.093	105	2039	0.441	ug/l	88
78) 1,2,4-Trimethylbenzene	11.470	105	2147	0.460	ug/l	100
79) sec-Butylbenzene	11.643	105	3296	0.417	ug/l	95
81) p-Isopropyltoluene	11.798	119	2436	0.492	ug/l	96
84) n-Butylbenzene	12.212	91	6754	0.608	ug/l	92
87) 1,2,4-Trichlorobenzene	13.839	180	5278	0.446	ug/l	96
89) Naphthalene	14.087	128	21121	1.125	ug/l	96
90) 1,2,3-Trichlorobenzene	14.328	180	5722	0.532	ug/l	92

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

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