

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU081922\
 Data File : VU050385.D
 Acq On : 19 Aug 2022 23:14
 Operator : SY/MD
 Sample : N3734-12
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WS-TRIHALOMETHANES

Manual Integrations
 APPROVED

Reviewed By :Krupa
 Patel

Quant Time: Aug 20 06:30:07 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U081922DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Fri Aug 19 15:06:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Fluorobenzene	6.112	96	30192	1.000 ug/l	# 0.00	0.00

System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.636	95	9017	0.877 ug/l		0.00
Spiked Amount 1.000			Recovery	=	88.000%	
68) 1,2-Dichlorobenzene-d4	12.195	152	10663	0.877 ug/l		0.00
Spiked Amount 1.000			Recovery	=	88.000%	

Target Compounds					Qvalue	
14) Carbon Disulfide	2.790	76	9708	0.352 ug/l	# 95	
15) Methylene Chloride	3.044	84	9464	0.787 ug/l		92
27) Chloroform	5.089	83	245593	13.240 ug/l		99
31) Isopropyl Ether	4.002	45	6838	0.230 ug/l		92
43) Dibromomethane	6.919	93	2625	0.487 ug/l		95
55) Dibromochloromethane	8.809	129	271416	31.130 ug/l		100
67) Bromoform	10.292	173	113299	22.883 ug/l		100
73) n-propylbenzene	10.909	120	3513	0.869 ug/l		93
89) Naphthalene	14.092	128	29869	1.885 ug/l		99

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

09/12/2022
 Supervised By :Mahesh
 Dadoda

09/12/2022

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