

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU082222\
 Data File : VU050398.D
 Acq On : 22 Aug 2022 13:56
 Operator : SY/MD
 Sample : N3734-12DL 5X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WS-TRISALOMETHANESDL

Quant Time: Aug 23 08:26:58 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U081922DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Tue Aug 23 08:18:20 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Fluorobenzene	6.118	96	31721	1.000	ug/l	# 0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.639	95	8437	0.781	ug/l	0.00
Spiked Amount 1.000			Recovery	=	78.000%	
68) 1,2-Dichlorobenzene-d4	12.198	152	10272	0.805	ug/l	0.00
Spiked Amount 1.000			Recovery	=	80.000%	
Target Compounds						
						Qvalue
15) Methylene Chloride	3.061	84	2838	0.225	ug/l	85
27) Chloroform	5.086	83	52785	2.708	ug/l	99
44) Bromodichloromethane	7.138	83	109652	8.202	ug/l	99
55) Dibromochloromethane	8.826	129	54605	5.961	ug/l	99
67) Bromoform	10.298	173	23603	4.537	ug/l	99

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

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