

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU091123\
 Data File : VU055236.D
 Acq On : 11 Sep 2023 15:39
 Operator : MD/SY
 Sample : 04354-01
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :

Quant Time: Sep 21 05:14:53 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U091123DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Thu Sep 21 05:13:10 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Fluorobenzene	6.116	96	51191	1.000	ug/l	# 0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.637	95	17775	0.895	ug/l	0.00
Spiked Amount 1.000			Recovery	=	89.000%	
68) 1,2-Dichlorobenzene-d4	12.196	152	18029	0.931	ug/l	0.00
Spiked Amount 1.000			Recovery	=	93.000%	
Target Compounds						
						Qvalue
3) Chloromethane	1.518	50	4504	0.256	ug/l	# 86
13) Acetone	2.628	43	3398	0.863	ug/l	92
24) tert-Butyl Alcohol	3.222	59	378	0.506	ug/l	# 1
27) Chloroform	5.087	83	54874	1.928	ug/l	93
46) t-1,4-Dichloro-2-butene	10.634	75	9133	3.465	ug/l	# 18
55) Dibromochloromethane	8.807	129	8558	0.727	ug/l	95
67) Bromoform	10.290	173	6806	1.147	ug/l	# 96
71) 1,2,3-Trichloropropane	10.634	75	9133	0.961	ug/l	# 33
89) Naphthalene	14.093	128	12386	0.503	ug/l	# 94

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

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