

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU091123\
 Data File : VU055232.D
 Acq On : 11 Sep 2023 13:18
 Operator : MD/SY
 Sample : VU0911WBL01
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :

Quant Time: Sep 21 05:14:14 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	6.113	96	58915	1.000	ug/l	# 0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.634	95	20098	0.879	ug/l	0.00
Spiked Amount 1.000			Recovery	=	88.000%	
68) 1,2-Dichlorobenzene-d4	12.193	152	18839	0.845	ug/l	0.00
Spiked Amount 1.000			Recovery	=	85.000%	
Target Compounds						
						Qvalue
13) Acetone	2.640	43	1247	0.275	ug/l	# 84
24) tert-Butyl Alcohol	3.187	59	612	0.711	ug/l	# 1
45) 4-Methyl-2-Pentanone	7.808	43	2106	0.260	ug/l	# 90
46) t-1,4-Dichloro-2-butene	10.637	75	9888	3.260	ug/l	# 18
54) 2-Hexanone	8.695	43	2368	0.304	ug/l	# 35
71) 1,2,3-Trichloropropane	10.637	75	9888	0.904	ug/l	# 34
87) 1,2,4-Trichlorobenzene	13.843	180	5215	0.261	ug/l	96
89) Naphthalene	14.090	128	18234	0.644	ug/l	99
90) 1,2,3-Trichlorobenzene	14.331	180	5163	0.298	ug/l	98

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

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