

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_R\DATA\VR111918\
Data File : VR026226.D
Acq On : 19 Nov 2018 17:51
Operator : SY/MD
Sample : J5997-09
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_R
ClientSampleId :
E3YA7

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR110918WMA.M
Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	8.979	1182	1200	1209	rBV5	1793664	7846623	2.50%	1.163%
2	10.062	1366	1378	1387	rBV5	55522047	313523965	100.00%	46.456%
3	10.452	1438	1442	1452	rVB3	3684000	7041537	2.25%	1.043%
4	10.817	1498	1502	1512	rVB	1901362	3488609	1.11%	0.517%
5	11.407	1591	1599	1610	rBV	51230310	204571153	65.25%	30.312%
6	11.516	1611	1617	1628	rVB2	26319546	90322325	28.81%	13.384%
7	11.839	1663	1670	1678	rBV	13248717	29625265	9.45%	4.390%
8	12.137	1714	1719	1738	rVB	1996322	4996042	1.59%	0.740%
9	12.477	1769	1775	1783	rBV	1572156	3498943	1.12%	0.518%
10	12.775	1819	1824	1834	rVB	1651972	3368889	1.07%	0.499%
11	12.922	1843	1848	1863	rBV	3184637	6593264	2.10%	0.977%

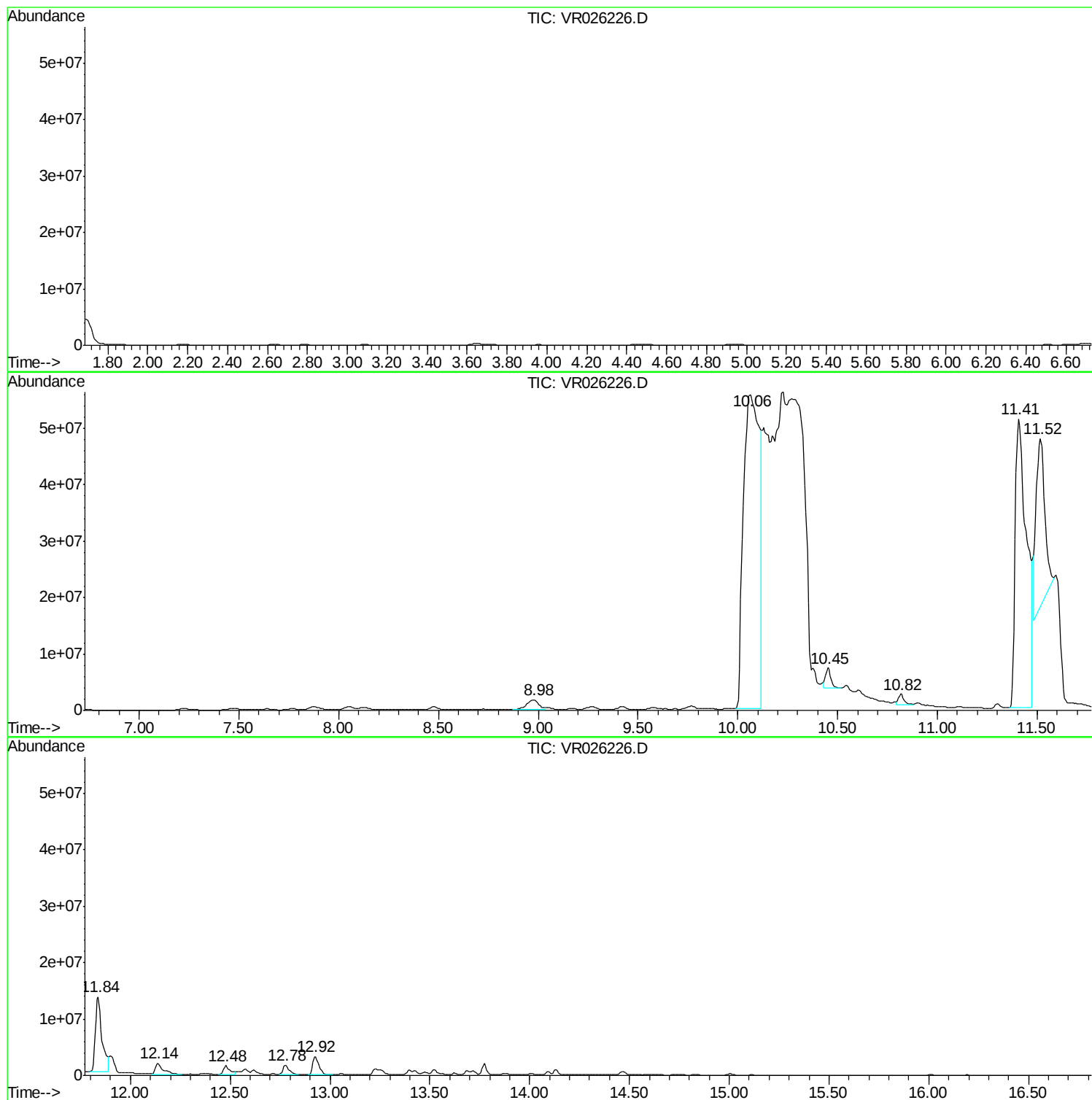
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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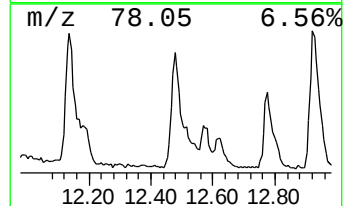
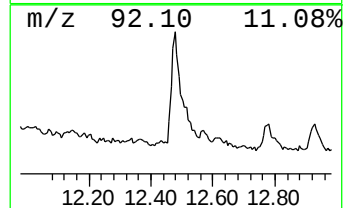
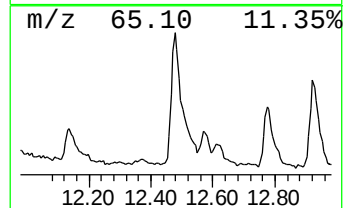
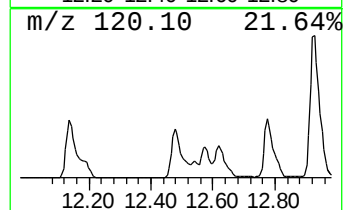
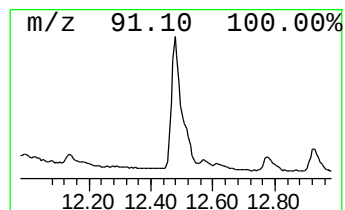
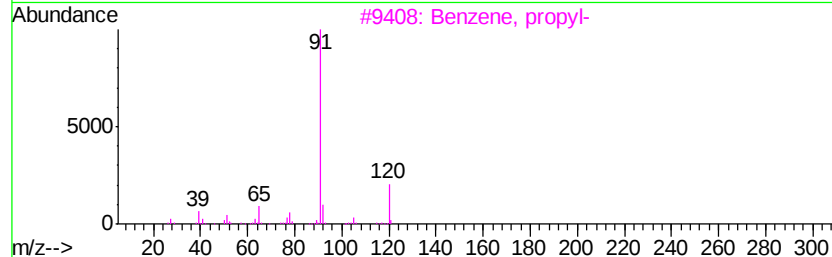
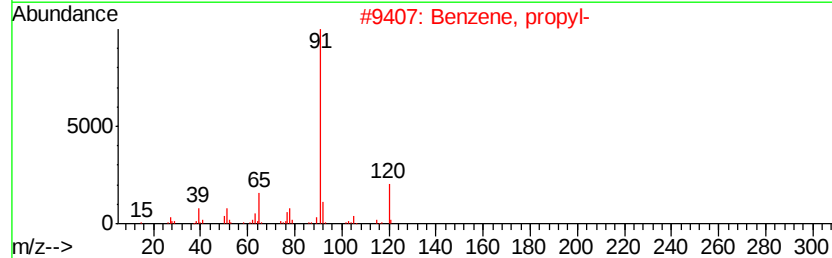
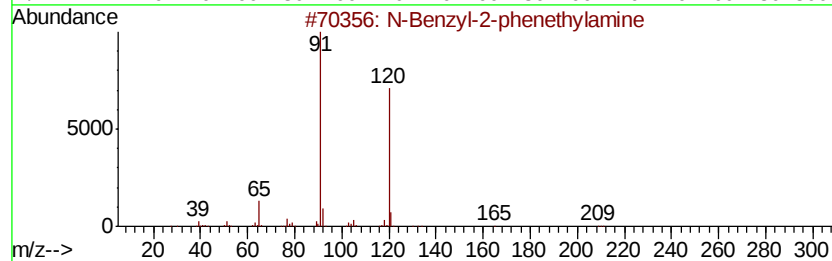
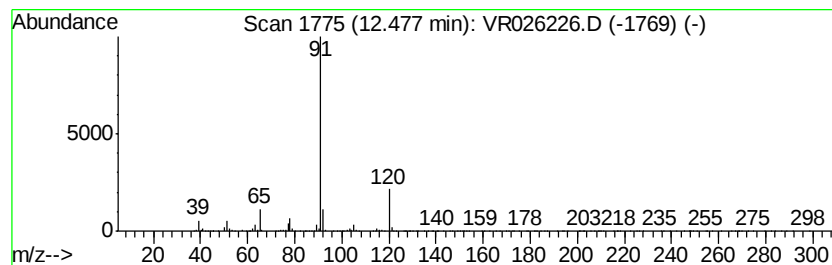
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 N-Benzyl-2-phenethylamine Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.48	2.65 ug/L	3498940	1,4-Dichlorobenzene-d4		13.23	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	91
2		Benzene, propyl-	120	C9H12	000103-65-1	91
3		Benzene, propyl-	120	C9H12	000103-65-1	90
4		Benzene, propyl-	120	C9H12	000103-65-1	87
5		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	87



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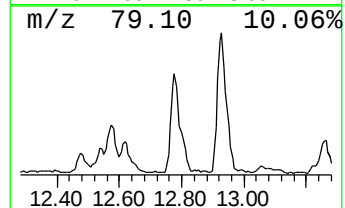
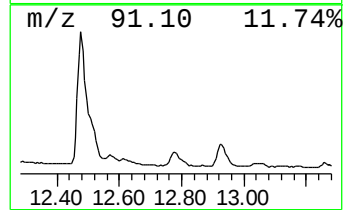
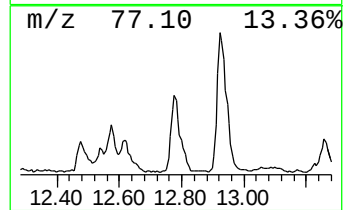
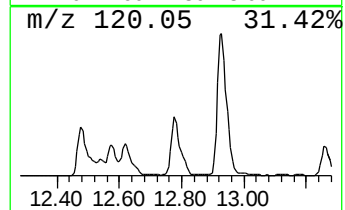
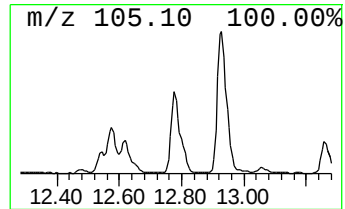
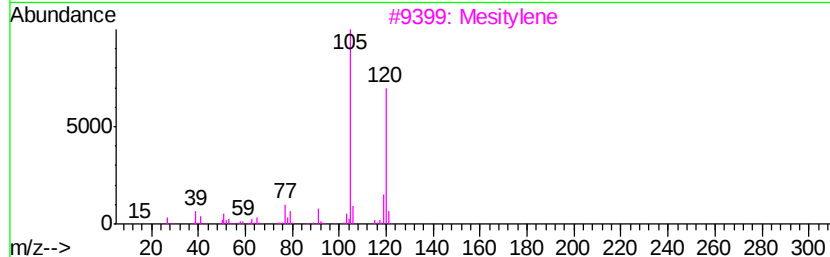
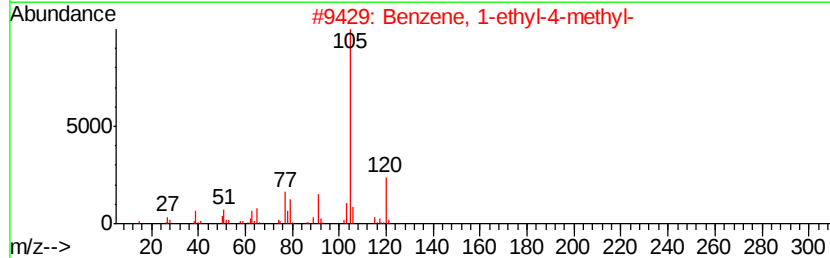
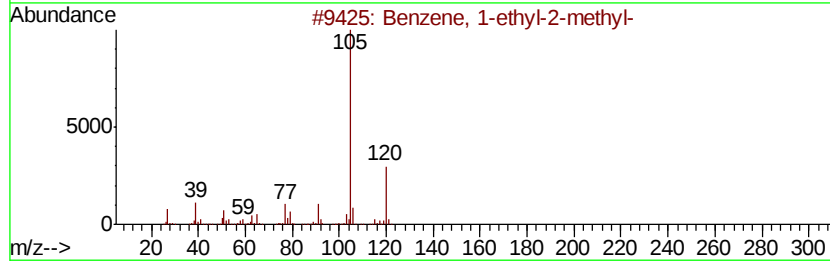
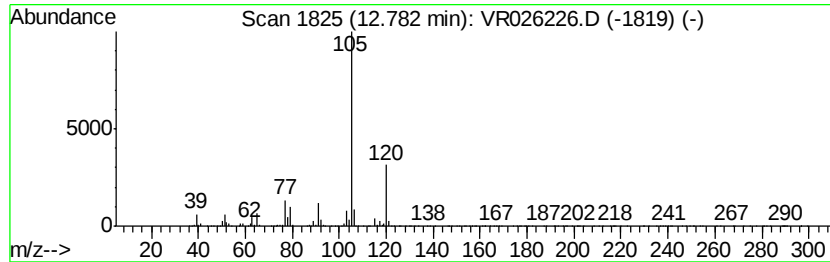
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Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.78	2.55 ug/L	3368890	1,4-Dichlorobenzene-d4	13.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3		Mesitylene	120	C9H12	000108-67-8	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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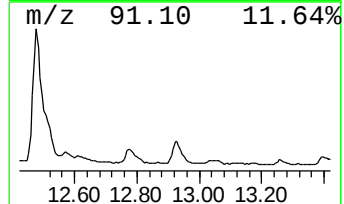
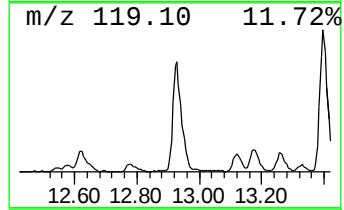
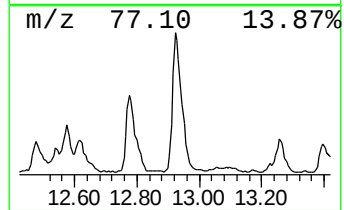
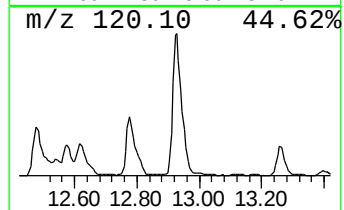
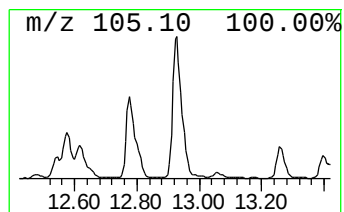
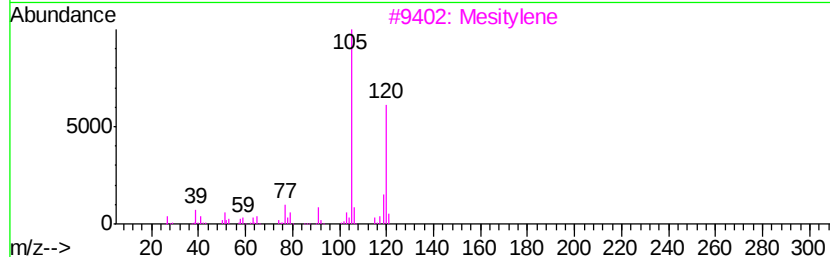
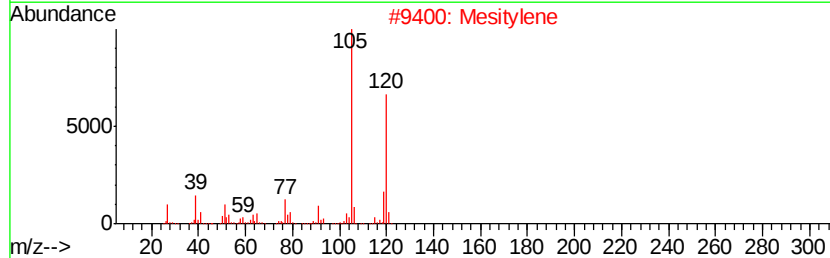
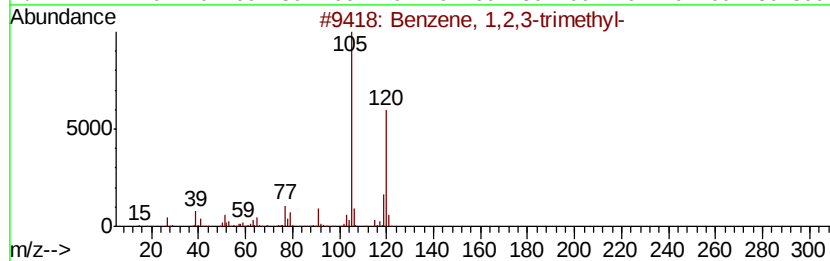
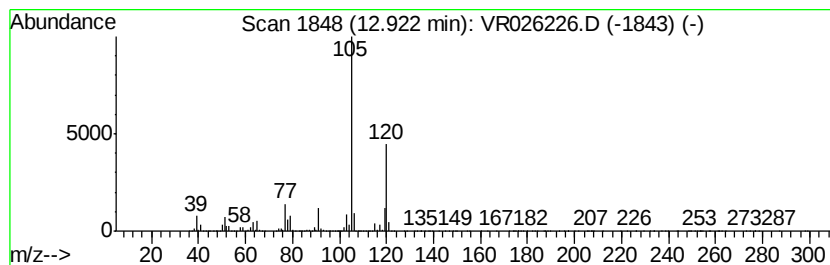
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Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.92	5.00 ug/L	6593260	1,4-Dichlorobenzene-d4	13.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Mesitylene	120	C9H12	000108-67-8	94
3		Mesitylene	120	C9H12	000108-67-8	94
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
N-Benzyl-2-phenet...	12.48	2.6	ug/L	3498940	3	13.23	6593260	5.0
Benzene, 1-ethyl...	12.78	2.5	ug/L	3368890	3	13.23	6593260	5.0
Benzene, 1,2,3-tr...	12.92	5.0	ug/L	6593260	3	13.23	6593260	5.0