

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111117\
 Data File : BF100460.D
 Acq On : 12 Nov 2017 1:18
 Operator : SJ/JU
 Sample : I6369-05
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 110617-JETT-F-D-B

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:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.210	17	21	26	rVB	69971	90147	2.29%	0.371%
2	5.122	512	516	524	rBV	303077	309966	7.87%	1.276%
3	5.516	579	583	586	rBV	1594058	1425776	36.18%	5.868%
4	6.516	749	753	757	rBV2	1107773	1019695	25.88%	4.196%
5	6.669	775	779	782	rBV	2549932	2312065	58.67%	9.515%
6	6.904	815	819	822	rBV	1021264	868931	22.05%	3.576%
7	7.063	841	846	849	rVB	3164622	2977543	75.56%	12.254%
8	7.469	909	915	918	rBV	2321166	2478108	62.88%	10.198%
9	8.181	1033	1036	1040	rBV	1165161	1048556	26.61%	4.315%
10	8.733	1127	1130	1134	rBV	118958	93268	2.37%	0.384%
11	9.263	1214	1220	1223	rBV	3795644	3940834	100.00%	16.218%
12	9.645	1282	1285	1290	rBV	140819	118599	3.01%	0.488%
13	9.939	1329	1335	1344	rVB	1115770	1010763	25.65%	4.160%
14	10.604	1445	1448	1451	rBV	116039	85774	2.18%	0.353%
15	10.645	1452	1455	1459	rBV	273014	247800	6.29%	1.020%
16	10.727	1464	1469	1472	rBV	1279637	1243427	31.55%	5.117%
17	11.427	1584	1588	1591	rBV	863452	783904	19.89%	3.226%
18	13.010	1853	1857	1861	rBV	2922796	2856192	72.48%	11.754%
19	13.898	2004	2008	2015	rBV3	37976	48270	1.22%	0.199%
20	14.063	2032	2036	2044	rBV	888310	780628	19.81%	3.213%
21	15.545	2283	2288	2298	rBV	400031	558520	14.17%	2.299%

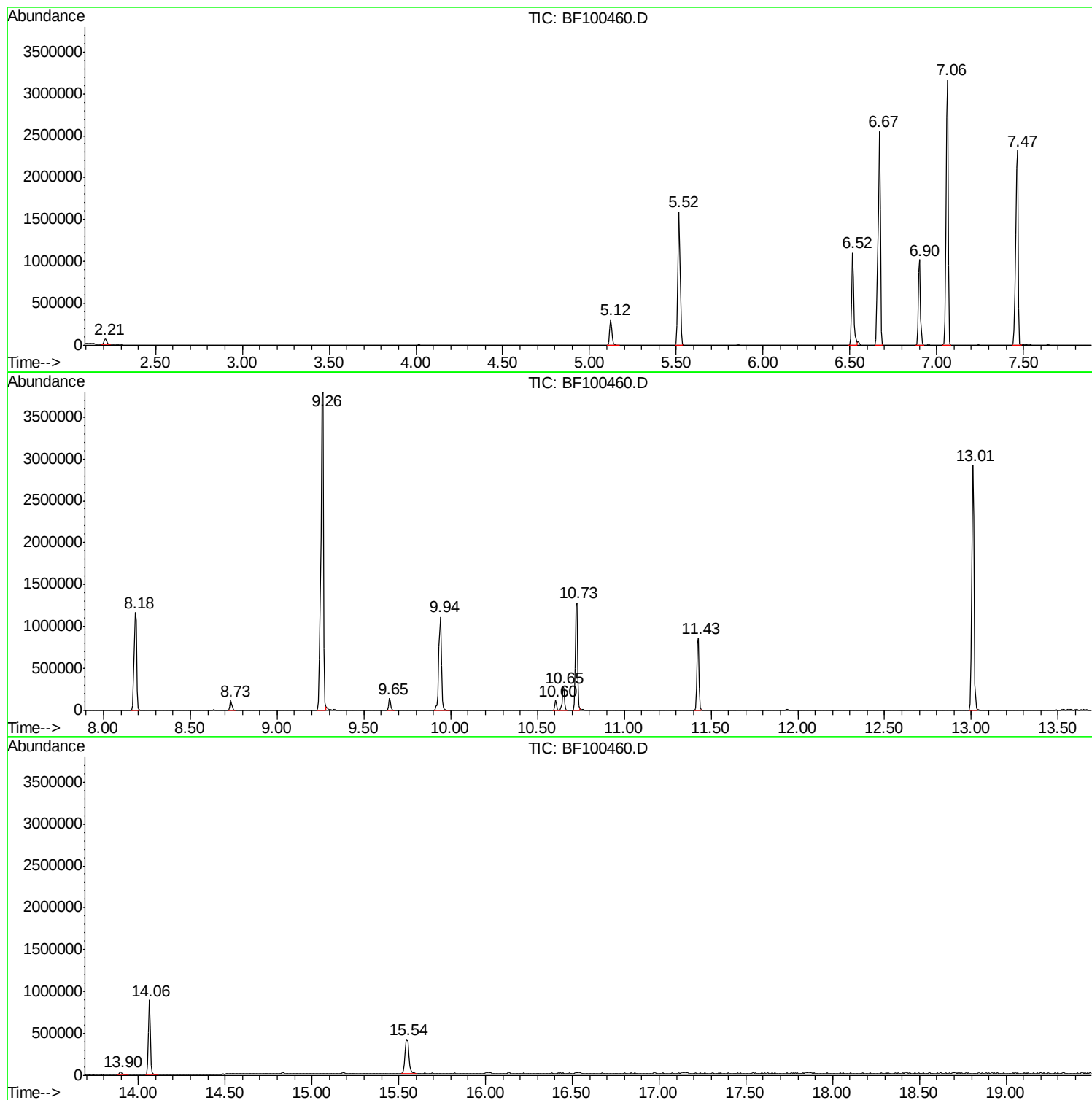
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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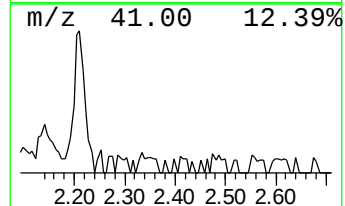
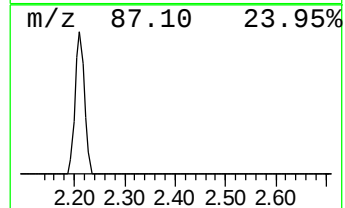
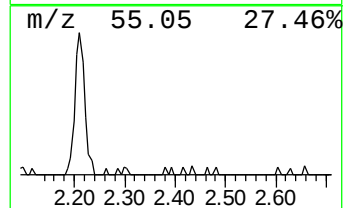
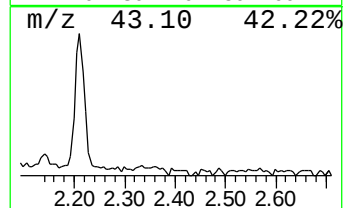
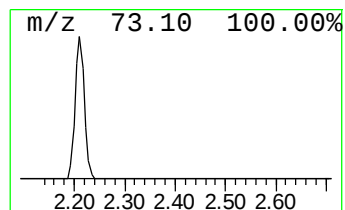
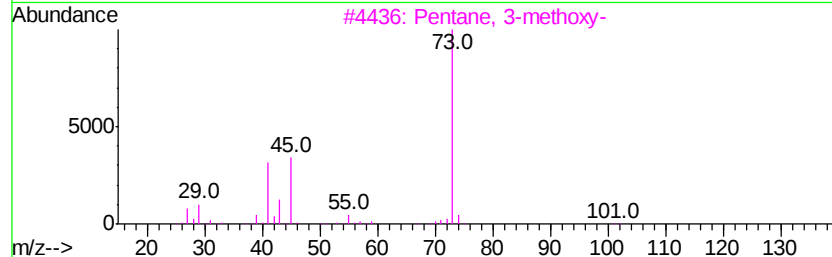
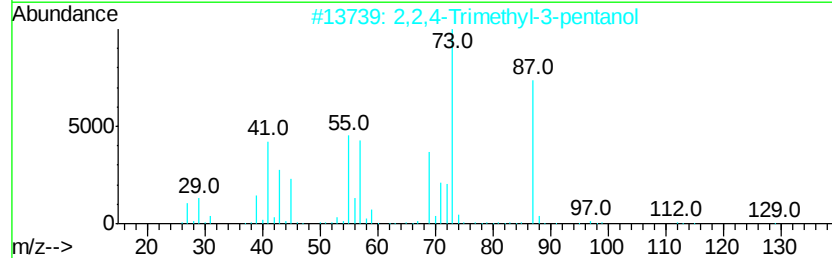
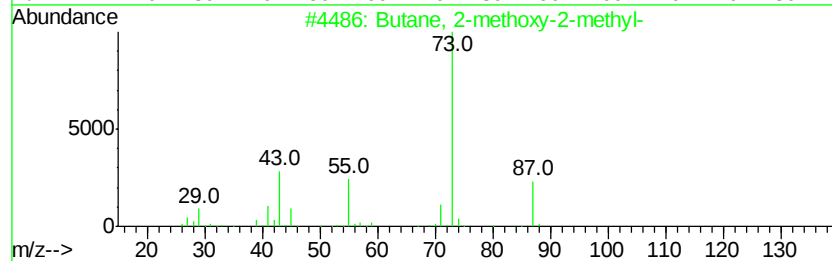
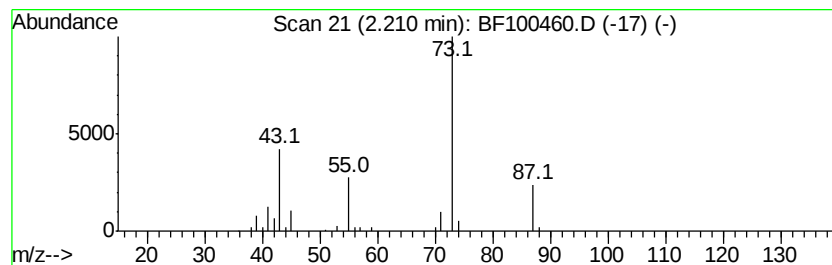
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.21	2.07 ng	90147	1,4-Dichlorobenzene-d4	6.90	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4	N-Ethylformamide	73	C3H7NO	000627-45-2	9
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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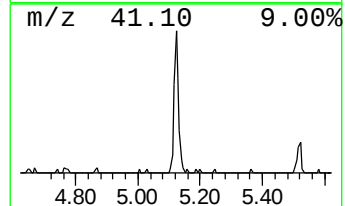
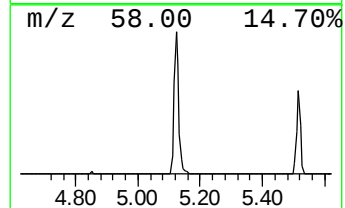
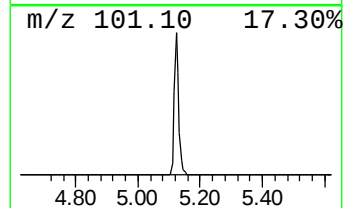
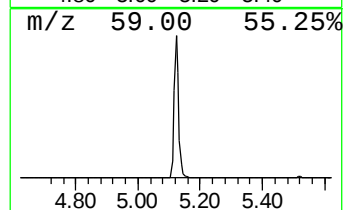
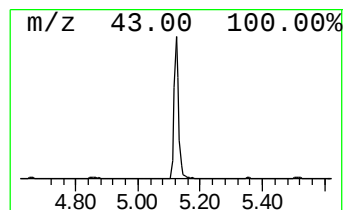
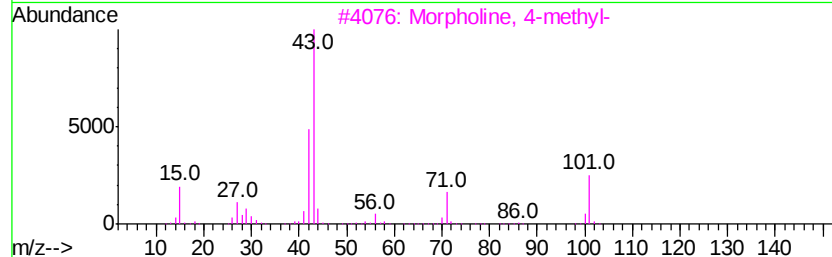
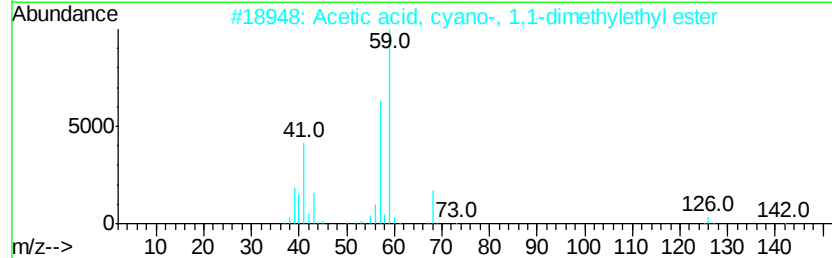
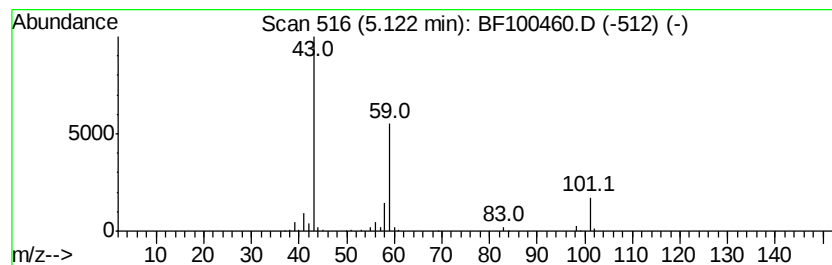
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	7.13 ng	309966	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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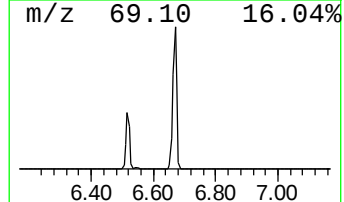
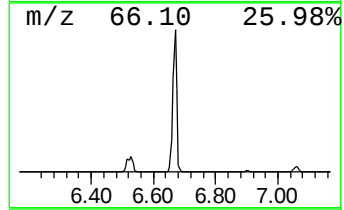
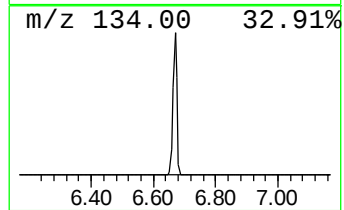
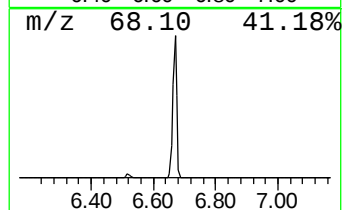
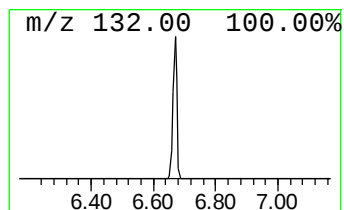
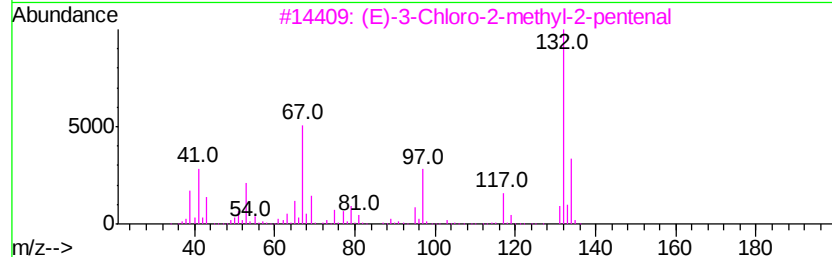
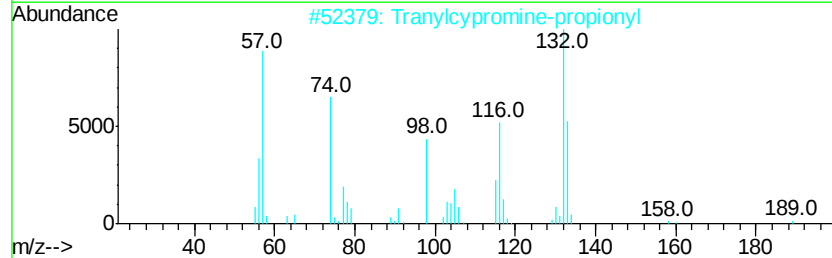
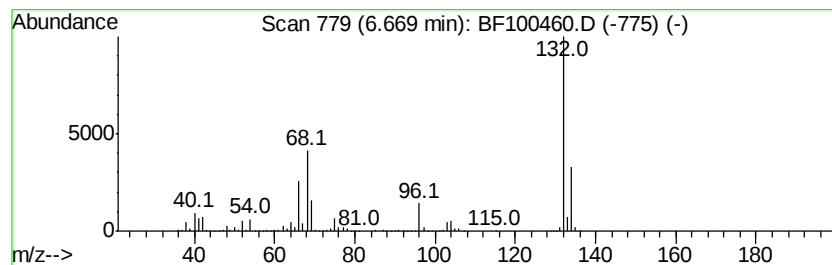
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Peak Number 3 unknown6.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	53.22 ng	2312070	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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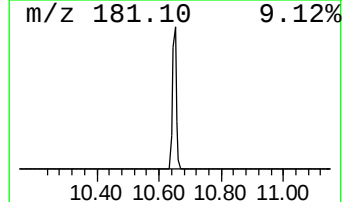
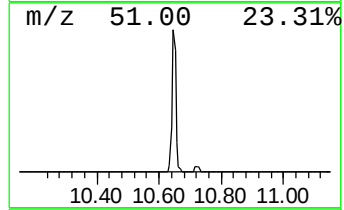
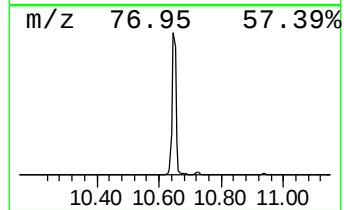
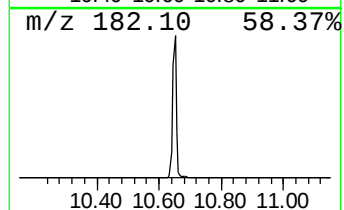
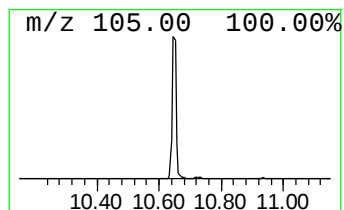
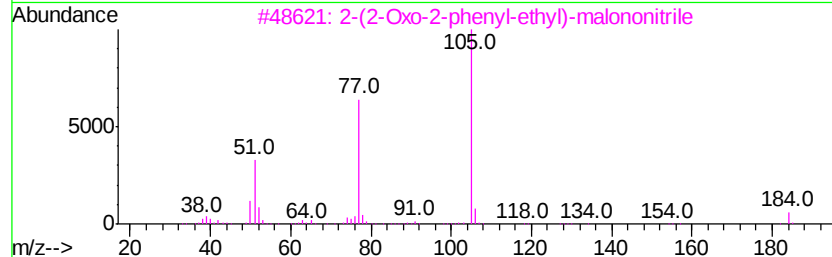
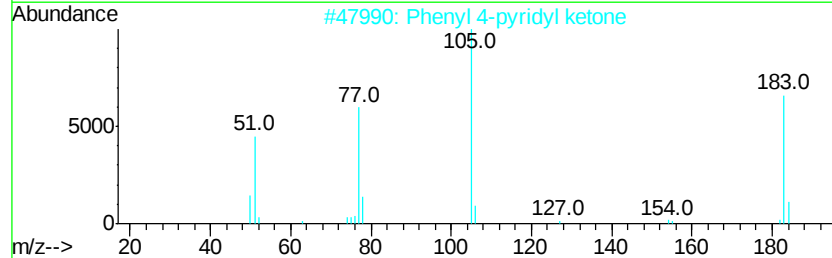
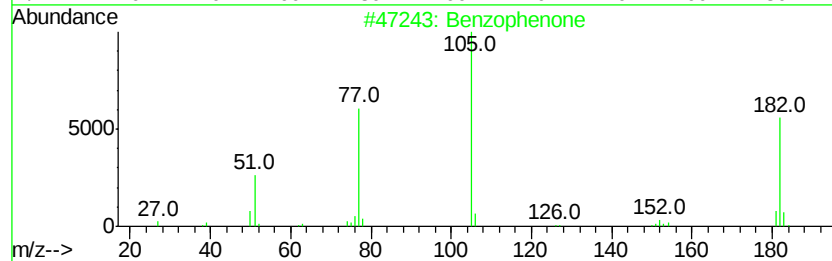
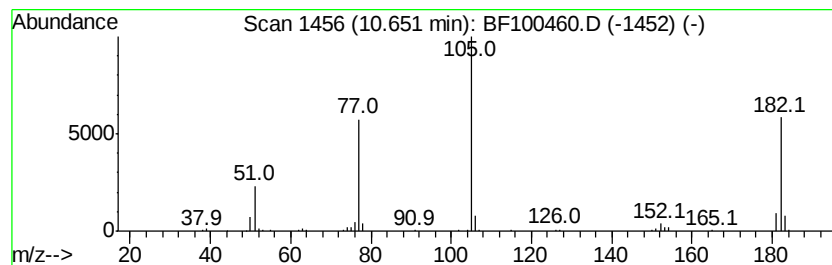
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Peak Number 5 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	4.90 ng	247800	Acenaphthene-d10	9.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzophenone	182	C13H10O	000119-61-9	96
2		Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	52
3		2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	47
4		Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47
5		Vinyl benzoate	148	C9H8O2	000769-78-8	47



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.21	2.1	ng	90147	1	6.90	868931	20.0
2-Pentanone, 4-hy...	5.12	7.1	ng	309966	1	6.90	868931	20.0
unknown6.67	6.67	53.2	ng	2312070	1	6.90	868931	20.0
Benzophenone	10.65	4.9	ng	247800	3	9.94	1010760	20.0