

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031015\
 Data File : BF077667.D
 Acq On : 10 Mar 2015 18:09
 Operator : TP/IZ
 Sample : G1466-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DCW-WW-087-3915

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: CENTROID

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
 Title : A

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.470	48	51	54	rVB	303956	522819	25.91%	3.654%
2	1.630	63	65	68	rVV	316678	447756	22.19%	3.129%
3	1.676	68	69	73	rVV	74967	129991	6.44%	0.908%
4	1.733	73	74	78	rVV	24904	33252	1.65%	0.232%
5	1.801	78	80	105	rVB	842054	1628954	80.73%	11.384%
6	4.944	354	355	361	rBV	47774	69469	3.44%	0.486%
7	5.470	399	401	404	rBV	543217	563157	27.91%	3.936%
8	6.750	510	513	515	rBV	246348	331095	16.41%	2.314%
9	6.887	523	525	528	rBV	887404	1171521	58.06%	8.188%
10	7.185	548	551	553	rBV	317625	293547	14.55%	2.052%
11	7.368	565	567	570	rVB	1365004	1236046	61.25%	8.639%
12	7.870	609	611	614	rBV	590186	790236	39.16%	5.523%
13	8.762	686	689	691	rBV	345762	345752	17.13%	2.416%
14	9.471	749	751	753	rBV	31198	30143	1.49%	0.211%
15	10.111	804	807	809	rBV	1366633	1751612	86.80%	12.242%
16	10.614	849	851	853	rVB	27573	27937	1.38%	0.195%
17	10.934	877	879	881	rVB	299090	390259	19.34%	2.727%
18	11.825	955	957	960	rBV	76113	89190	4.42%	0.623%
19	11.905	962	964	967	rBV	1050308	1009412	50.02%	7.055%
20	12.762	1037	1039	1042	rBV	385938	425346	21.08%	2.973%
21	14.763	1211	1214	1217	rBV	2041192	2017888	100.00%	14.103%
22	15.940	1314	1317	1321	rBV2	16673	24423	1.21%	0.171%
23	16.054	1324	1327	1330	rBV	463032	489628	24.26%	3.422%
24	17.826	1479	1482	1485	rBV	389355	489098	24.24%	3.418%

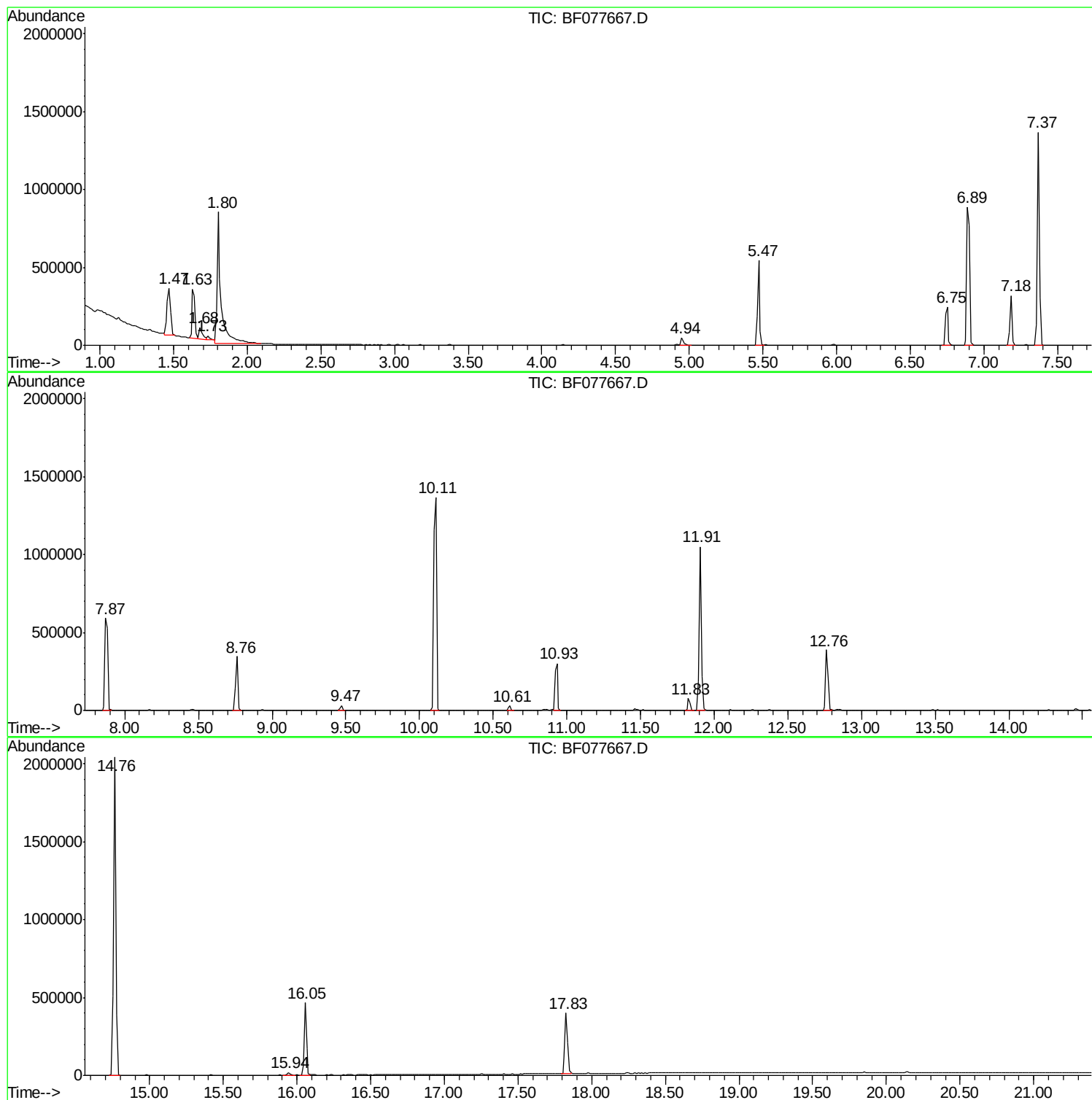
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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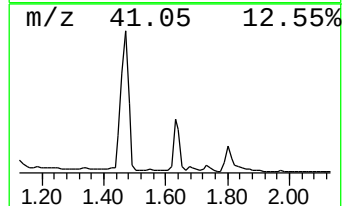
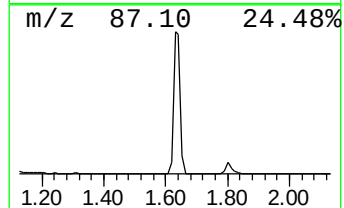
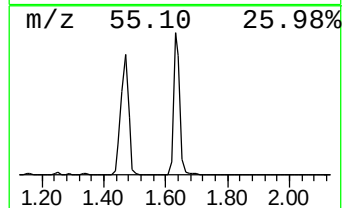
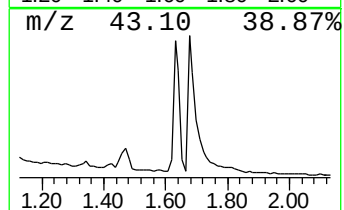
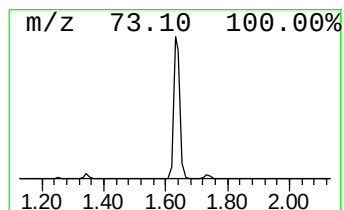
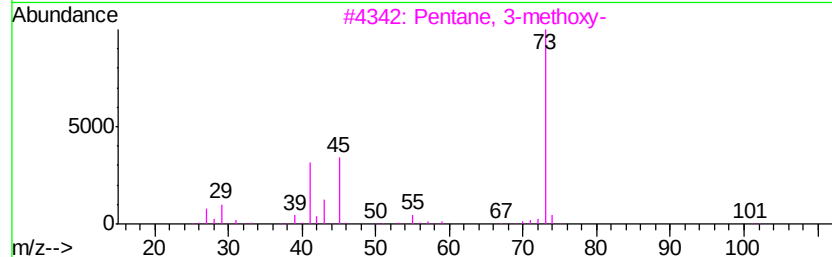
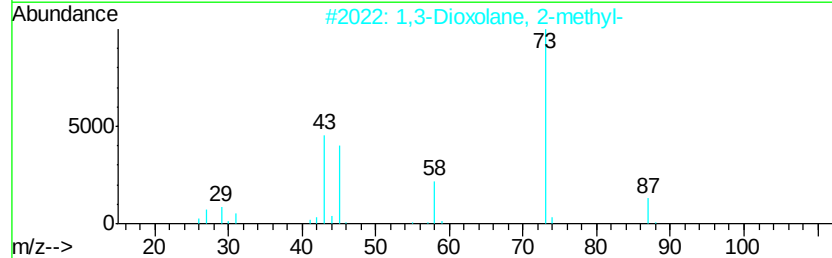
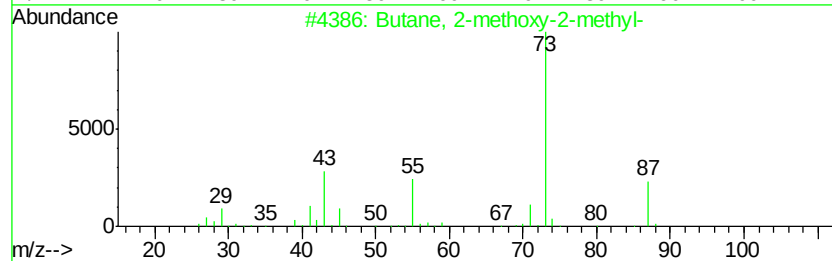
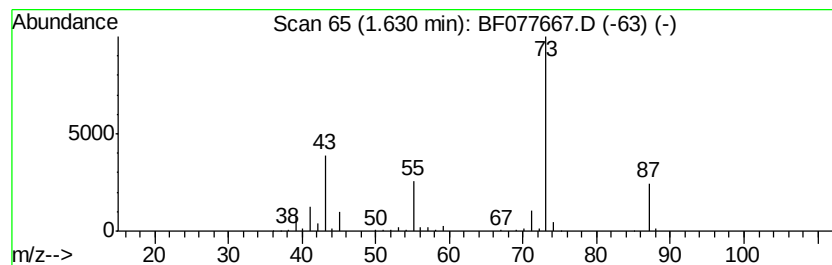
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF021915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.63	30.51 ng	447756	1,4-Dichlorobenzene-d4	7.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	10
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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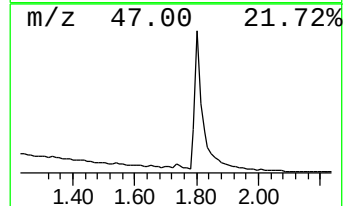
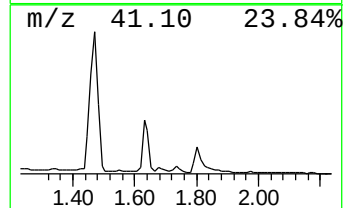
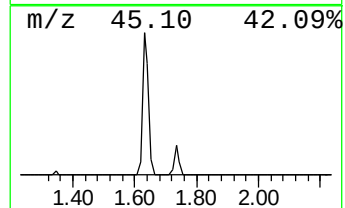
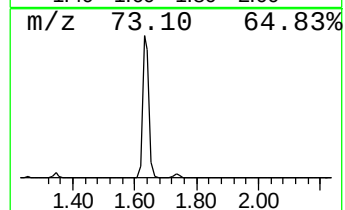
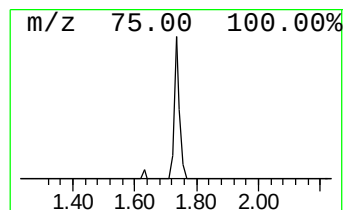
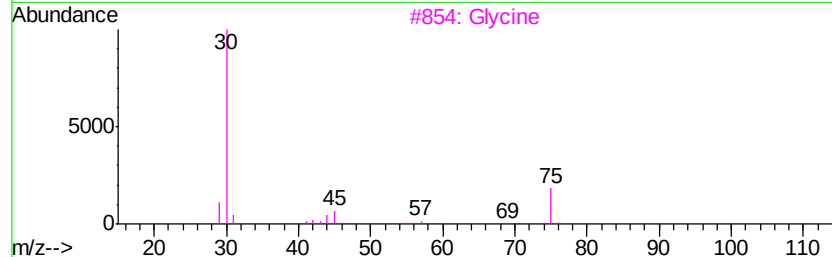
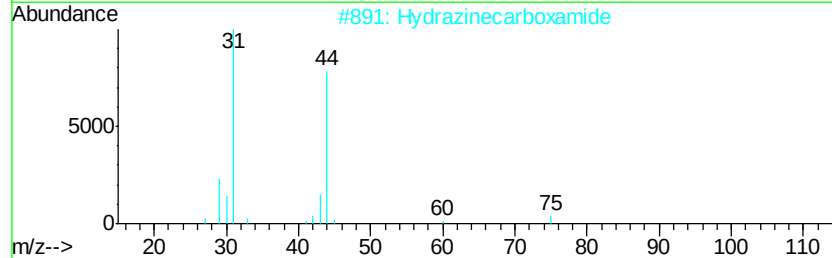
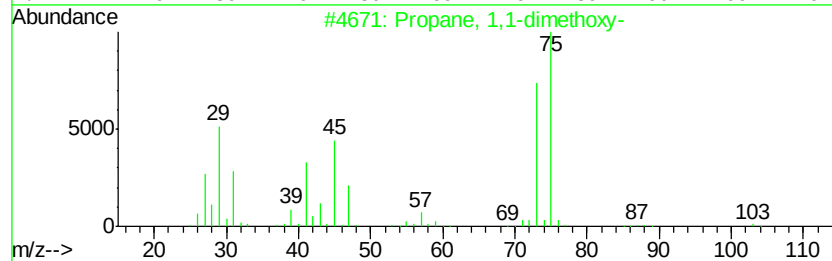
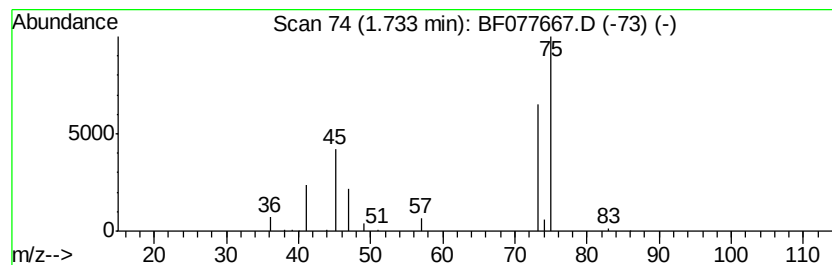
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Propane, 1,1-dimethoxy- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.73	2.27 ng	33252	1,4-Dichlorobenzene-d4	7.18

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	78
2		Hydrazinecarboxamide	75	CH5N3O	000057-56-7	3
3		Glycine	75	C2H5NO2	000056-40-6	3
4		Ethanethioamide	75	C2H5NS	000062-55-5	2
5		1,3-Dioxolane, 2-(chloromethyl)-	122	C4H7ClO2	002568-30-1	2



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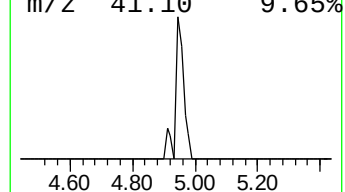
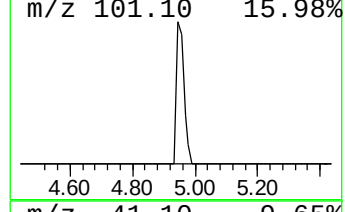
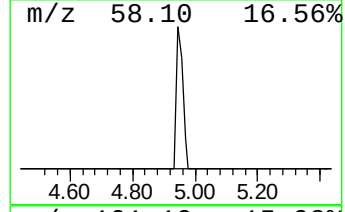
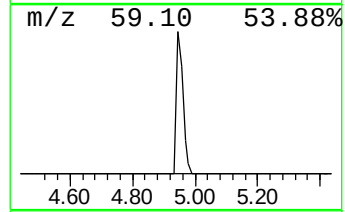
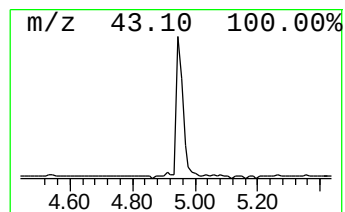
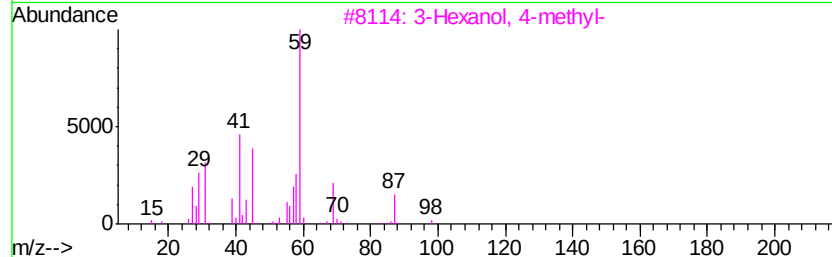
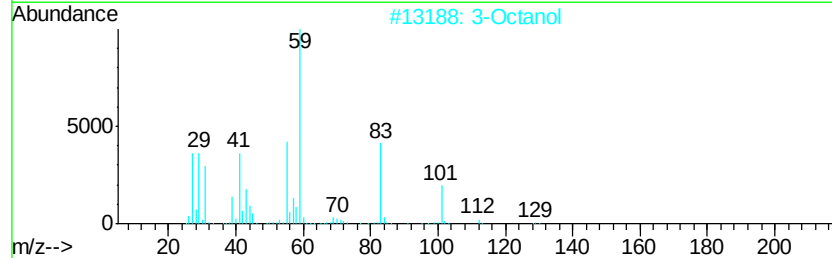
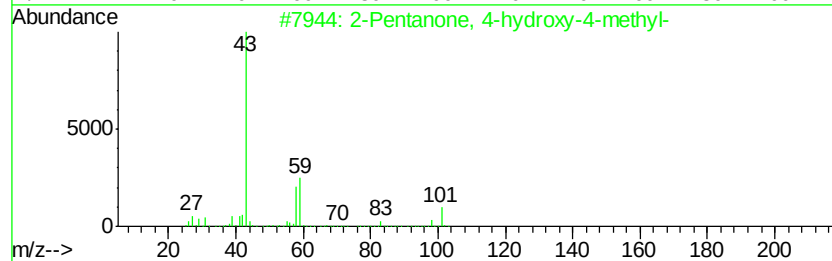
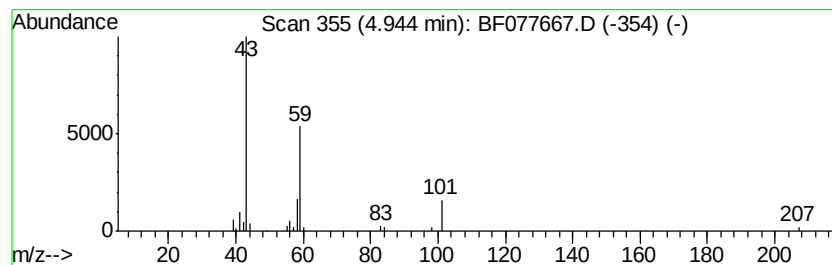
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.94	4.73 ng	69469	1,4-Dichlorobenzene-d4	7.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		3-Octanol	130	C8H18O	000589-98-0	36
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		Diacetamide	101	C4H7NO2	000625-77-4	9
5		Cycloserine	102	C3H6N2O2	000068-41-7	9



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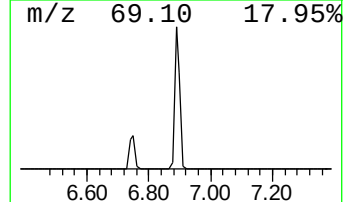
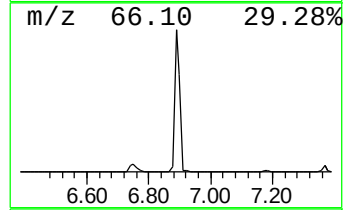
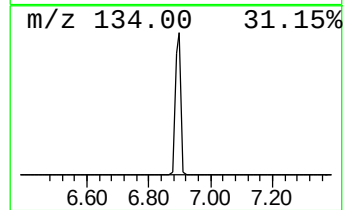
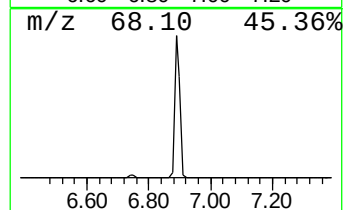
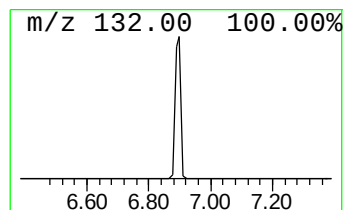
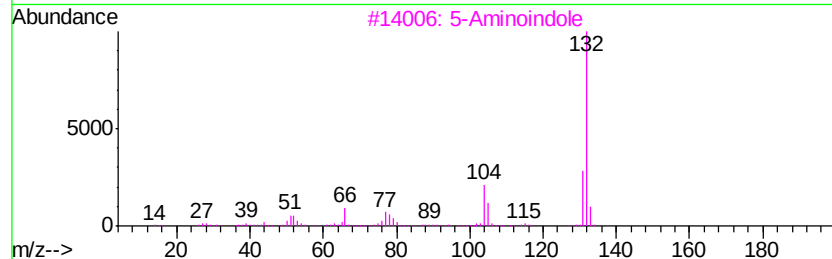
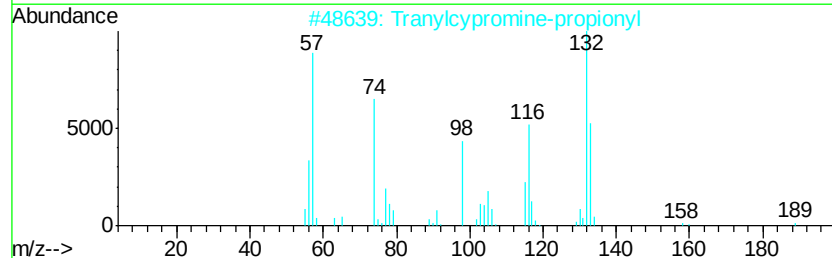
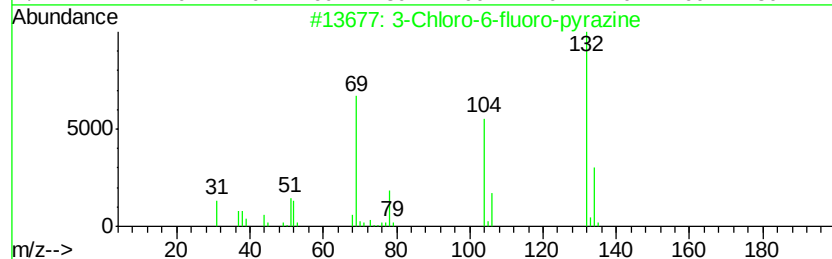
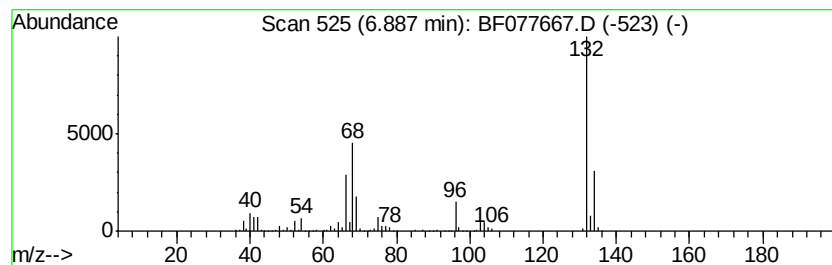
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Peak Number 7 unknown6.89 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.89	79.82 ng	1171520	1,4-Dichlorobenzene-d4	7.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
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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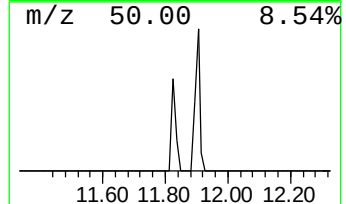
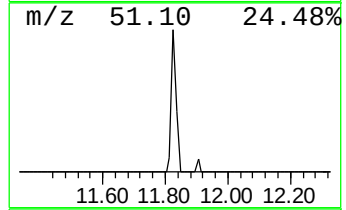
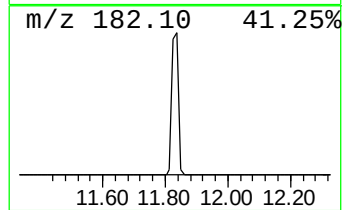
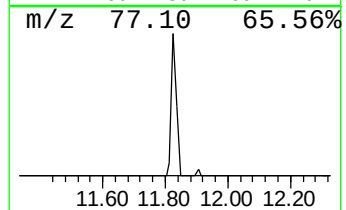
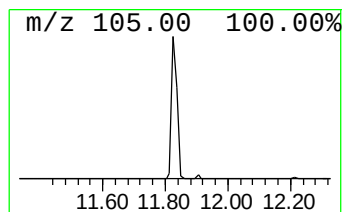
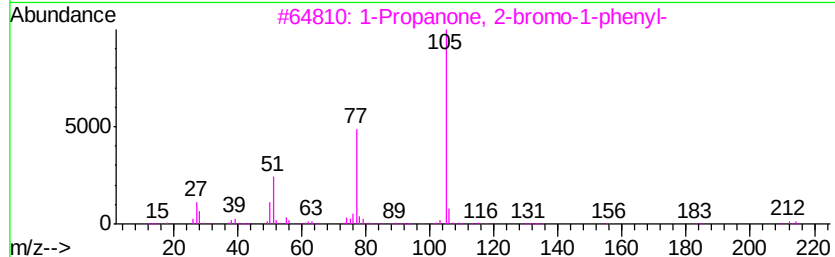
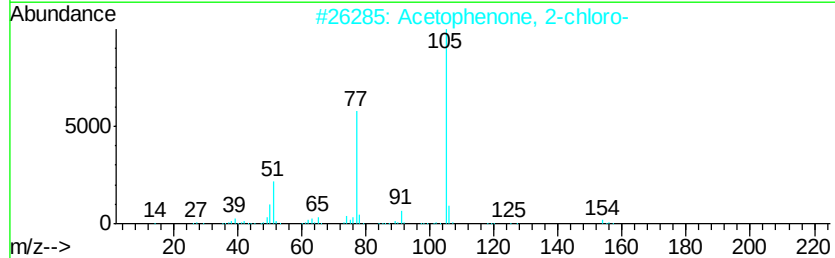
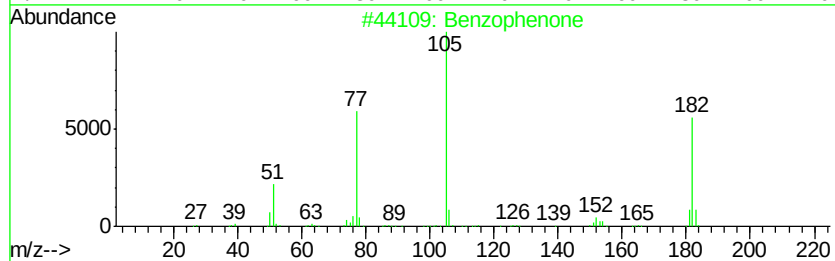
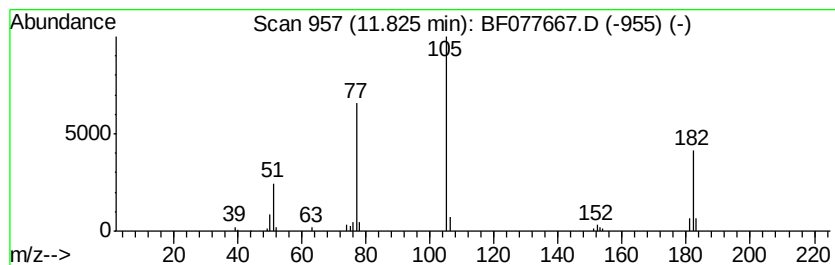
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Peak Number 9 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.83	4.57 ng	89190	Acenaphthene-d10	10.93
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzophenone	182 C13H10O	000119-61-9 96
2		Acetophenone, 2-chloro-	154 C8H7ClO	000532-27-4 49
3		1-Propanone, 2-bromo-1-phenyl-	212 C9H9BrO	002114-00-3 47
4		Benzenebutanoic acid, .gamma.-oxo-	178 C10H10O3	002051-95-8 47
5		4,5-Diphenyloxazolin-2(3H)-one	237 C15H11NO2	005014-83-5 47



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	1.63	30.5	ng	447756	1	7.18	293547	20.0
Propane, 1,1-dime...	1.73	2.3	ng	33252	1	7.18	293547	20.0
2-Pentanone, 4-hy...	4.94	4.7	ng	69469	1	7.18	293547	20.0
unknown6.89	6.89	79.8	ng	1171520	1	7.18	293547	20.0
Benzophenone	11.83	4.6	ng	89190	3	10.93	390259	20.0