

Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
 Data File : BF089496.D
 Acq On : 1 Aug 2016 5:31
 Operator : UM/SJ
 Sample : H4215-04
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-14-COMP

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.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	1.655	5	6	51	rVB	961975	2314347	100.00%	13.262%
2	4.581	259	262	273	rBV	237710	457198	19.75%	2.620%
3	5.061	302	304	320	rBV	913417	1237598	53.48%	7.092%
4	5.267	320	322	338	rVV	78030	161032	6.96%	0.923%
5	6.159	398	400	403	rBV	1043276	1251684	54.08%	7.172%
6	6.261	407	409	413	rVV	1318483	1323004	57.17%	7.581%
7	6.330	413	415	417	rVV	253493	280278	12.11%	1.606%
8	6.376	417	419	427	rVB	80432	146370	6.32%	0.839%
9	6.490	427	429	436	rVB	281834	291475	12.59%	1.670%
10	6.650	440	443	445	rBV	1042605	1043142	45.07%	5.977%
11	6.753	450	452	458	rVB	93654	124145	5.36%	0.711%
12	7.016	473	475	477	rBV	43602	42828	1.85%	0.245%
13	7.073	477	480	483	rVV	553758	825915	35.69%	4.733%
14	7.130	483	485	487	rVV	76035	127103	5.49%	0.728%
15	7.164	487	488	490	rVV	42268	58090	2.51%	0.333%
16	7.359	502	505	513	rVB	331340	376588	16.27%	2.158%
17	7.542	518	521	522	rBV	66205	82507	3.57%	0.473%
18	7.587	522	525	527	rVV	108241	146634	6.34%	0.840%
19	7.782	539	542	544	rBV	375633	378233	16.34%	2.167%
20	8.010	559	562	565	rBV	295013	356633	15.41%	2.044%
21	8.067	565	567	569	rBV	270705	368312	15.91%	2.110%
22	8.307	585	588	590	rVV3	10511	24282	1.05%	0.139%
23	8.479	602	603	608	rBV2	85915	132443	5.72%	0.759%
24	8.559	608	610	612	rBV	97444	94277	4.07%	0.540%
25	8.605	612	614	615	rBV	51217	44167	1.91%	0.253%
26	8.673	618	620	623	rBV2	64546	81245	3.51%	0.466%
27	8.776	628	629	632	rVB	73904	78729	3.40%	0.451%
28	8.856	634	636	639	rVB	996179	1215226	52.51%	6.963%
29	8.982	644	647	652	rBV2	101173	166009	7.17%	0.951%
30	9.142	660	661	663	rBV	25562	30281	1.31%	0.174%
31	9.256	669	671	673	rBV	192721	224519	9.70%	1.287%
32	9.473	688	690	693	rBV2	103389	205948	8.90%	1.180%
33	9.519	693	694	696	rVV	362369	387713	16.75%	2.222%
34	9.565	696	698	704	rVV	51871	104475	4.51%	0.599%

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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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35	9.702	707	710	712	rVV3	21642	30823	1.33%	0.177%
36	9.976	733	734	736	rVV	45177	36005	1.56%	0.206%
37	10.056	739	741	747	rVB6	11541	24378	1.05%	0.140%
38	10.193	749	753	756	rBV2	12891	31174	1.35%	0.179%
39	10.308	761	763	766	rVV	683652	660794	28.55%	3.786%
40	10.445	773	775	781	rVB	48015	60181	2.60%	0.345%
41	10.890	812	814	816	rBV	34684	32943	1.42%	0.189%
42	10.993	821	823	827	rBV	347366	326500	14.11%	1.871%
43	11.553	870	872	875	rBV	63310	64765	2.80%	0.371%
44	12.193	926	928	930	rBV	46551	44999	1.94%	0.258%
45	12.422	945	948	950	rBV	67416	71589	3.09%	0.410%
46	12.582	959	962	964	rBV	818523	942912	40.74%	5.403%
47	13.485	1039	1041	1050	rVB2	75751	112756	4.87%	0.646%
48	13.611	1050	1052	1058	rBV	294298	393598	17.01%	2.255%
49	14.125	1095	1097	1100	rVB3	15588	26866	1.16%	0.154%
50	14.605	1137	1139	1143	rBV	31110	59593	2.57%	0.341%
51	14.845	1158	1160	1163	rVB	23775	27456	1.19%	0.157%
52	14.902	1163	1165	1167	rVV	21380	27499	1.19%	0.158%
53	14.948	1167	1169	1179	rVB	207805	294300	12.72%	1.686%

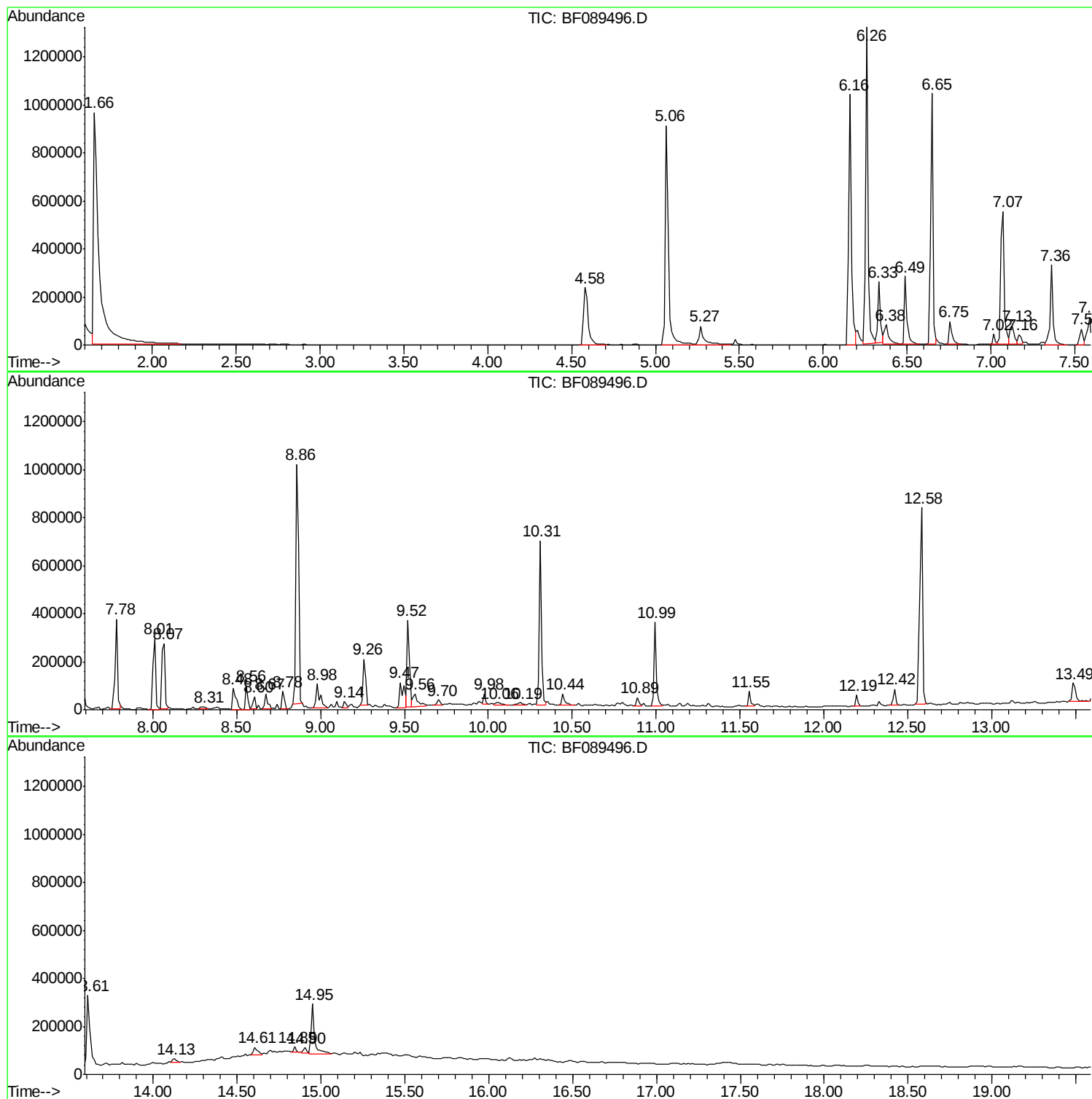
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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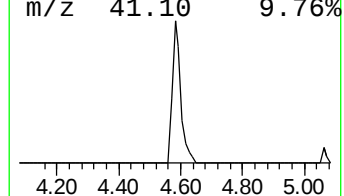
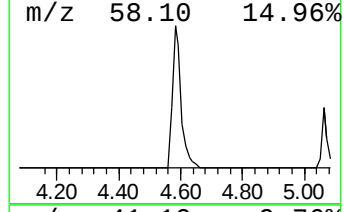
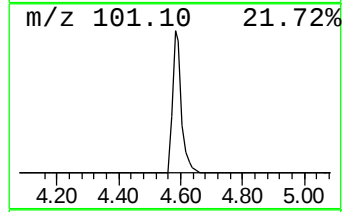
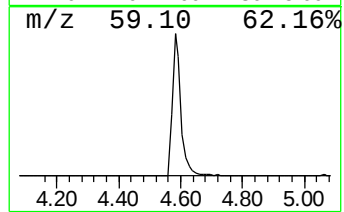
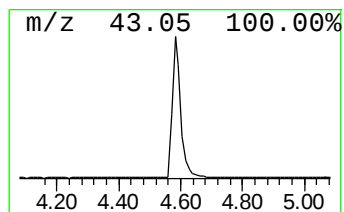
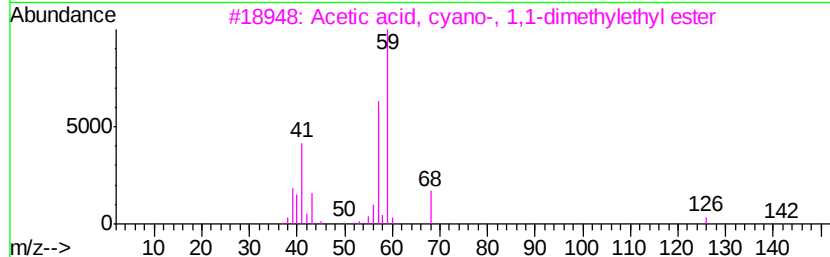
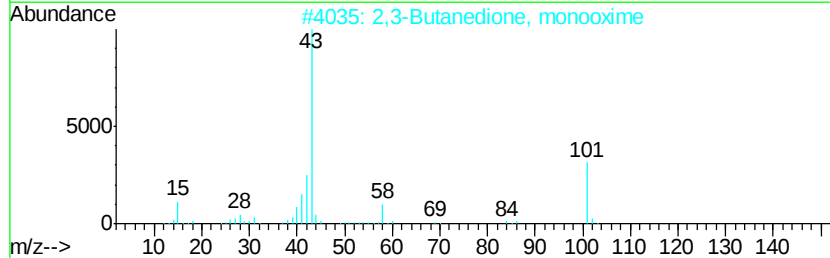
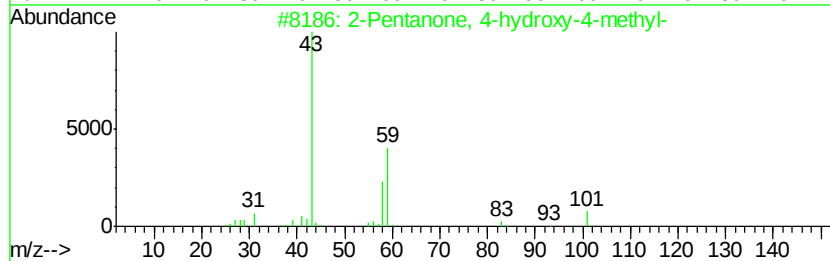
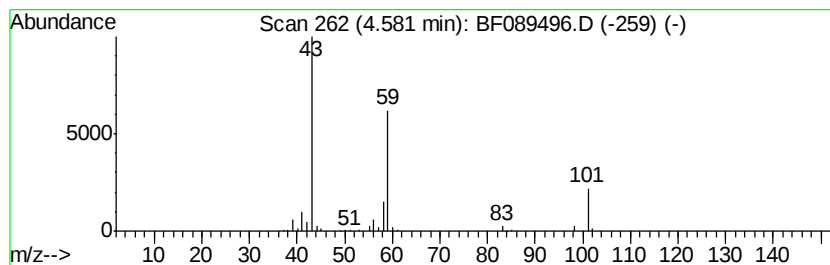
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.58	31.37 ng	457198	1,4-Dichlorobenzene-d4	6.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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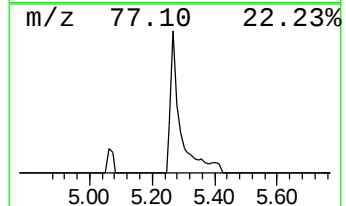
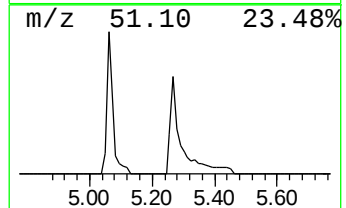
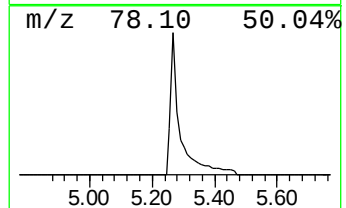
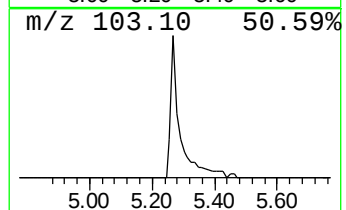
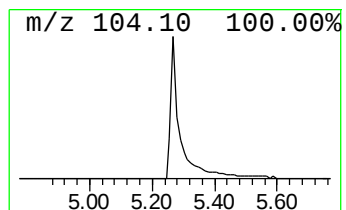
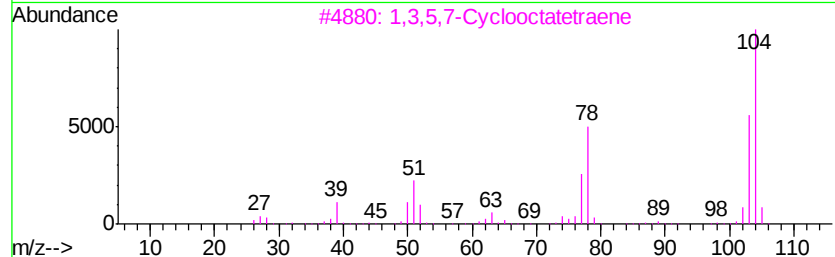
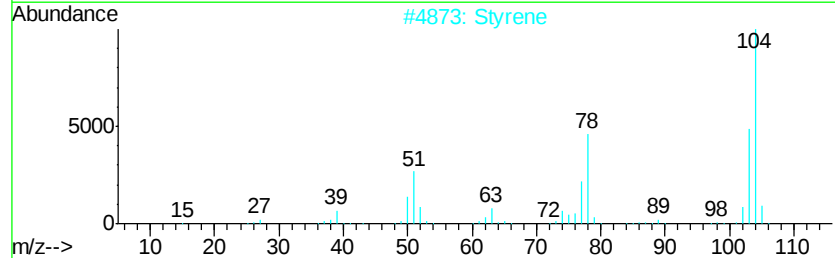
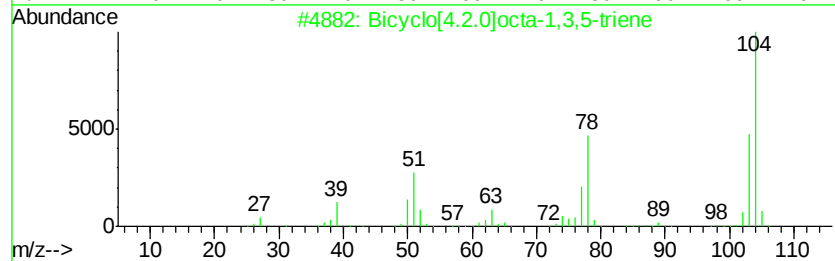
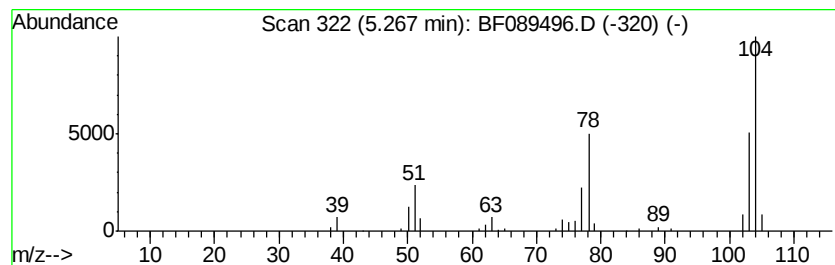
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Peak Number 3 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.27	11.05 ng	161032	1,4-Dichlorobenzene-d4	6.49

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Bicyclo[4.2.0]octa-1,3,5-triene	104	C8H8	000694-87-1	96
2		Styrene	104	C8H8	000100-42-5	96
3		1,3,5,7-Cyclooctatetraene	104	C8H8	000629-20-9	94
4		Butanedioic acid, phenyl-	194	C10H10O4	000635-51-8	72
5		Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	39



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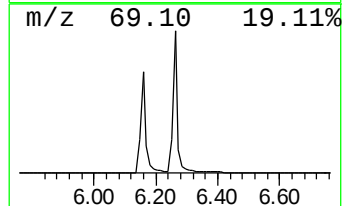
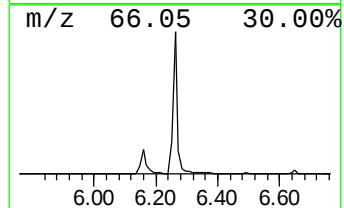
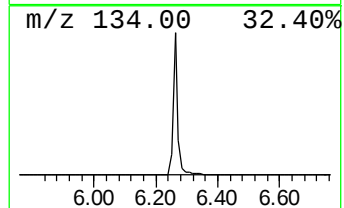
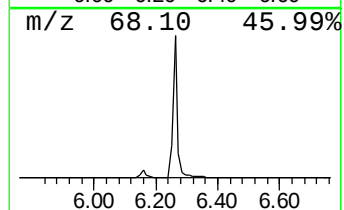
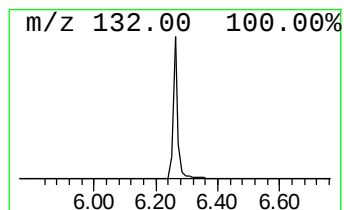
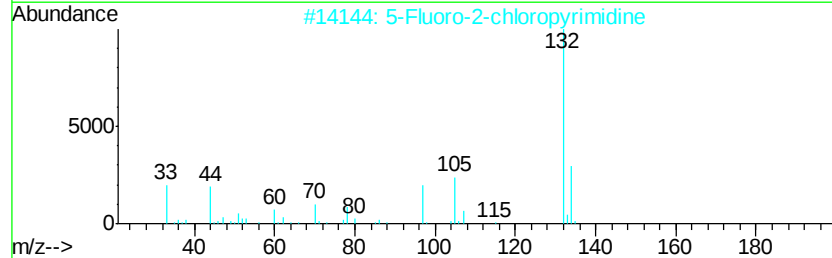
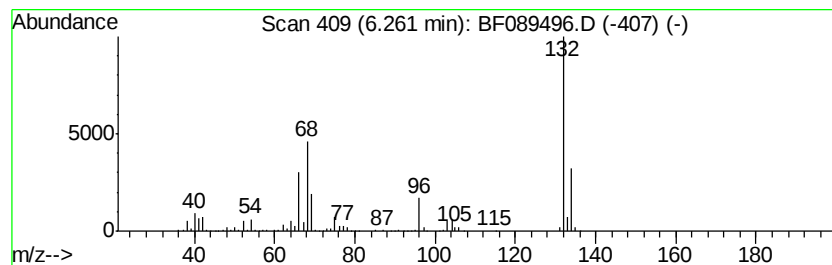
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown6.26 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.26	90.78 ng	1323000	1,4-Dichlorobenzene-d4	6.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		Xenon	132	Xe	007440-63-3	9



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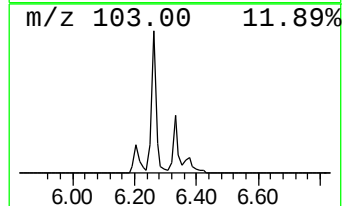
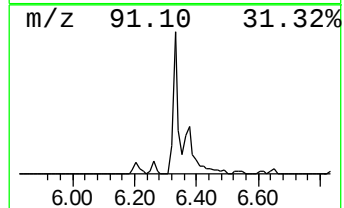
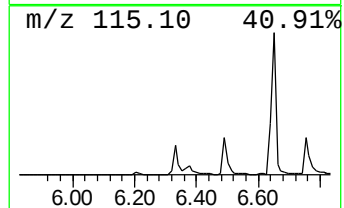
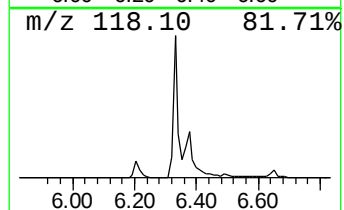
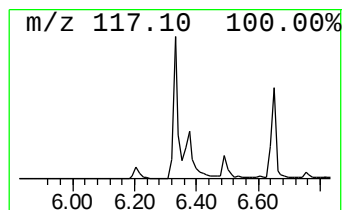
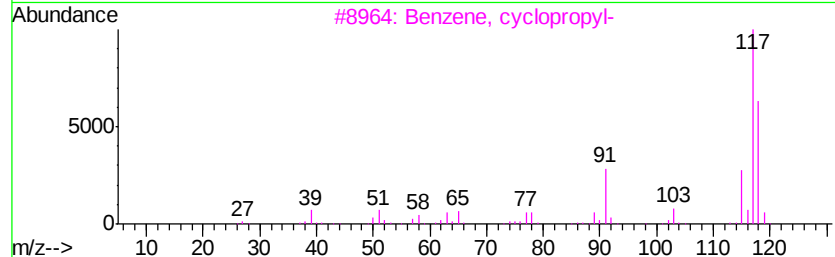
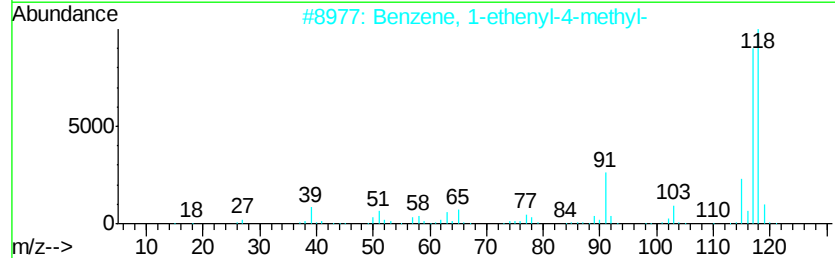
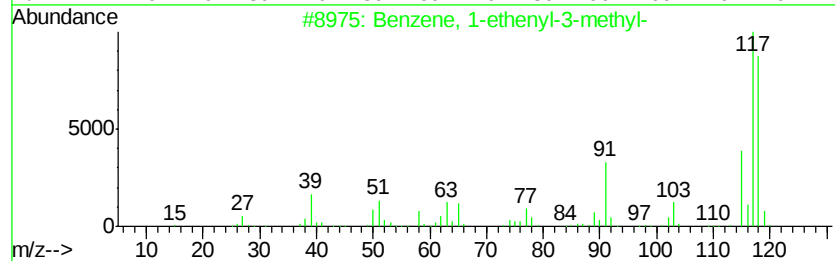
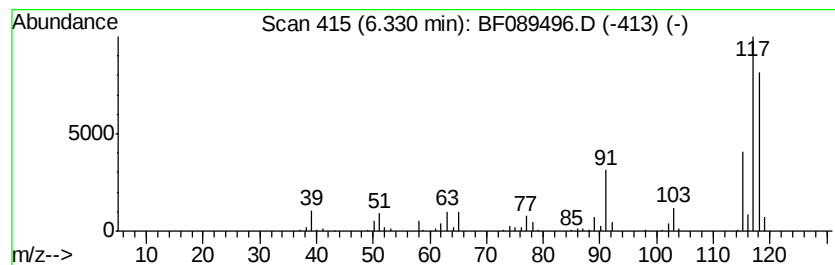
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethenyl-3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.33	19.23 ng	280278	1,4-Dichlorobenzene-d4	6.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	96
2		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	95
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	94
4		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	94
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	93



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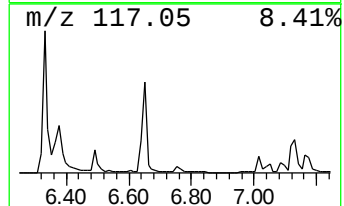
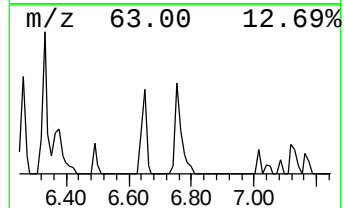
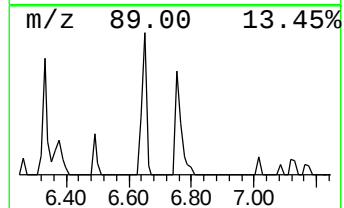
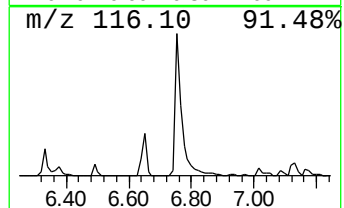
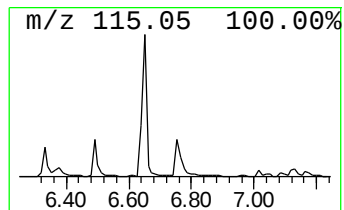
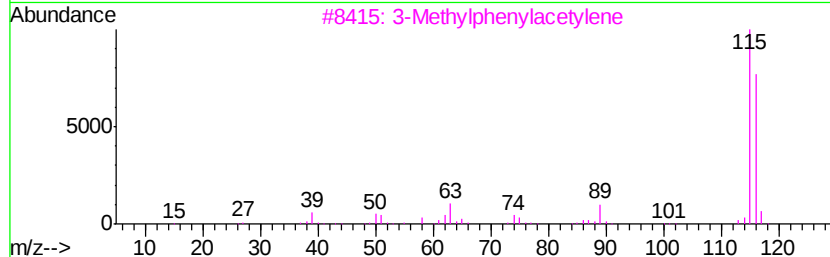
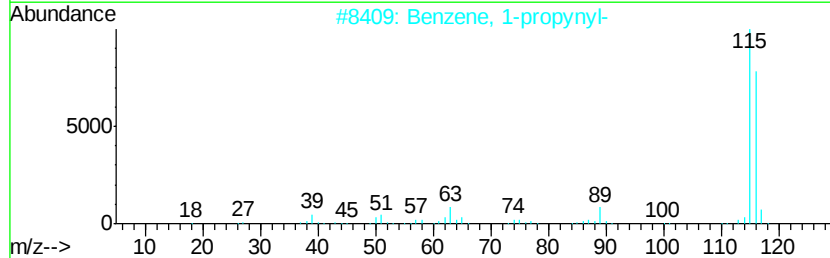
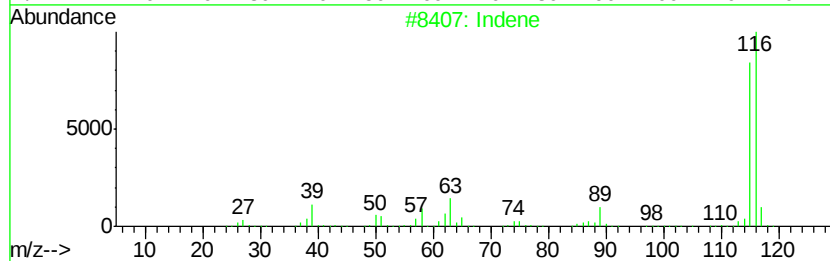
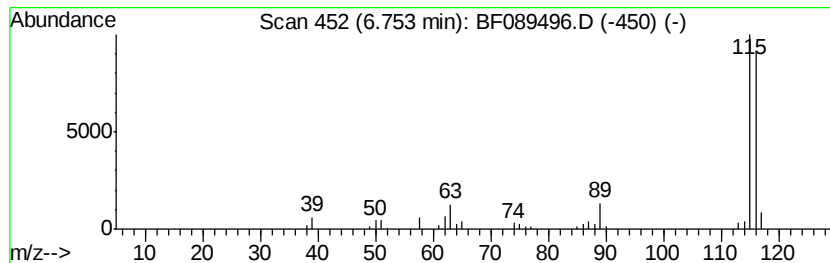
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Indene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.75	8.52 ng	124145	1,4-Dichlorobenzene-d4	6.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	95
2		Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
3		3-Methylphenylacetylene	116	C9H8	000766-82-5	91
4		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91
5		1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
Data File : BF089496.D
Acq On : 1 Aug 2016 5:31
Operator : UM/SJ
Sample : H4215-04
Misc :
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Instrument :
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ClientSampleId :
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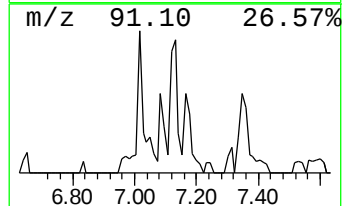
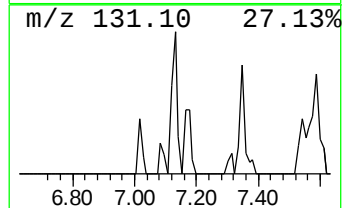
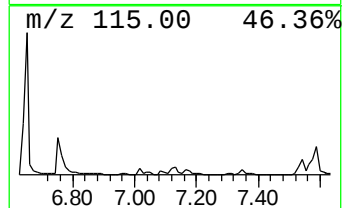
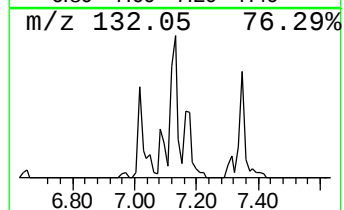
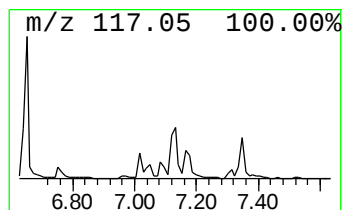
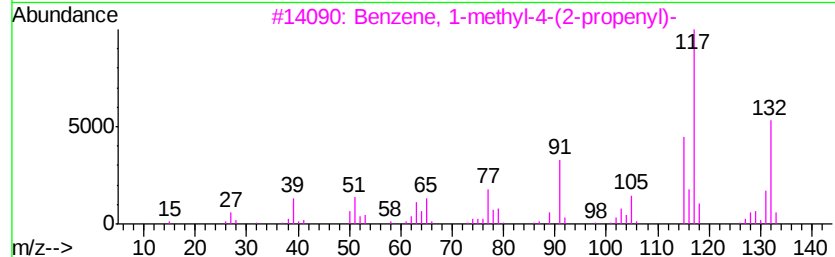
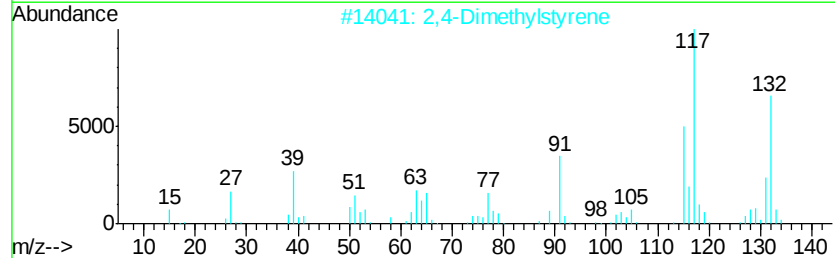
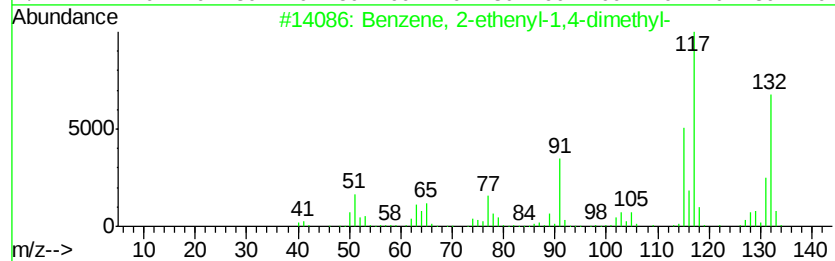
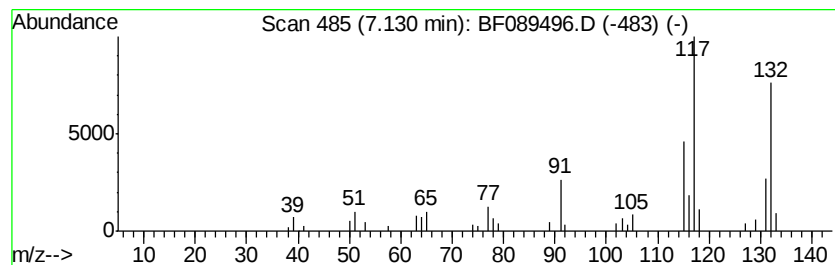
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	8.72 ng	127103	1,4-Dichlorobenzene-d4	6.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	97
2		2,4-Dimethylstyrene	132	C10H12	002234-20-0	96
3		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	95
4		(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	94
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF073116\
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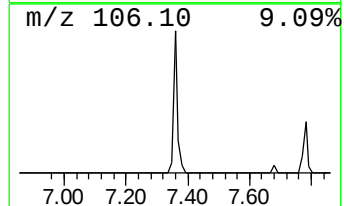
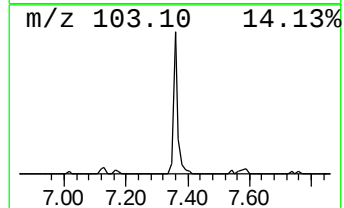
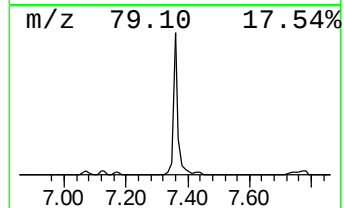
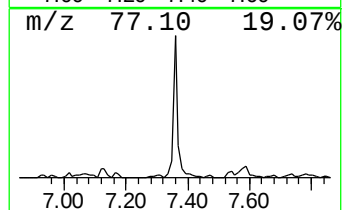
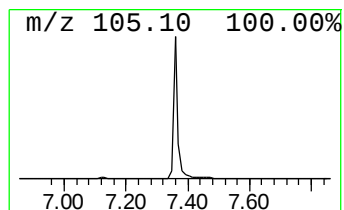
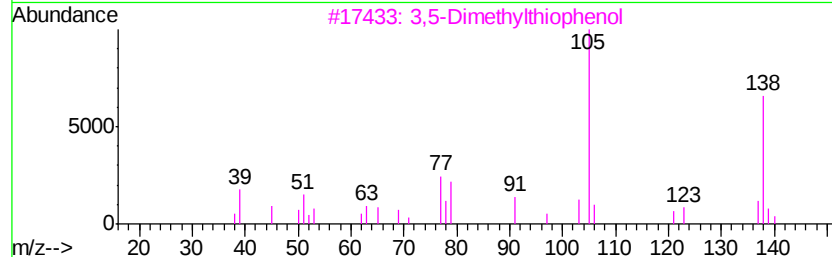
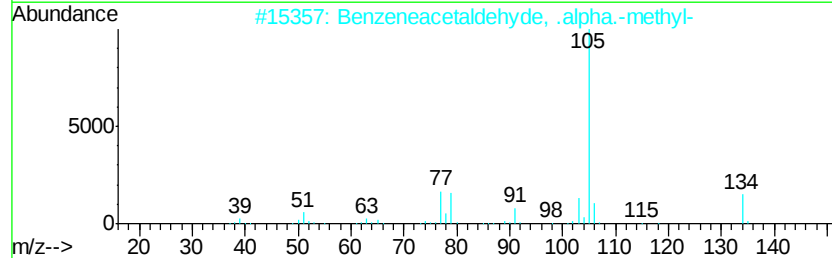
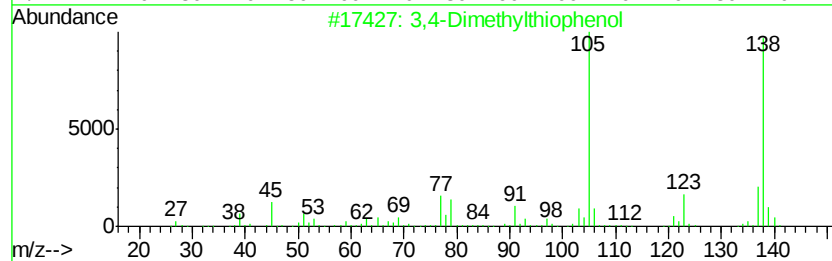
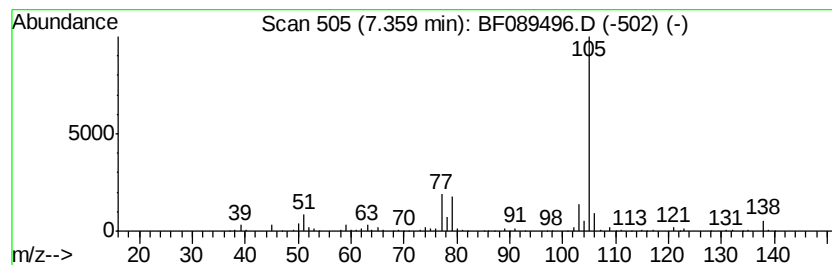
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 3,4-Dimethylthiophenol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.36	19.91 ng	376588	Napthalene-d8	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,4-Dimethylthiophenol	138	C8H10S	018800-53-8	78
2		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	72
3		3,5-Dimethylthiophenol	138	C8H10S	038360-81-5	72
4		3-Methylbenzyl mercaptan	138	C8H10S	025697-56-7	72
5		Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF073116\
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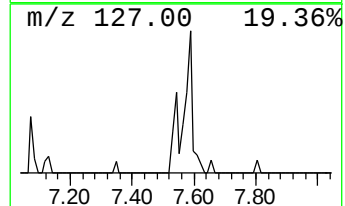
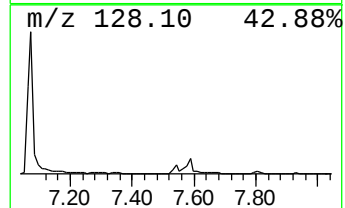
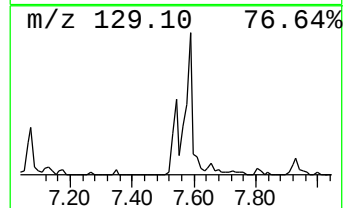
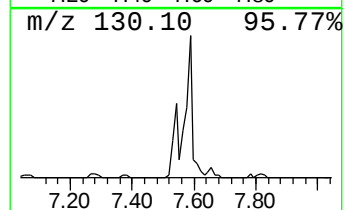
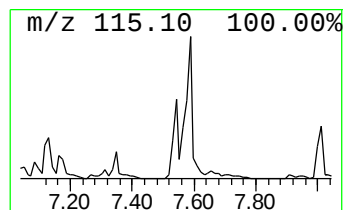
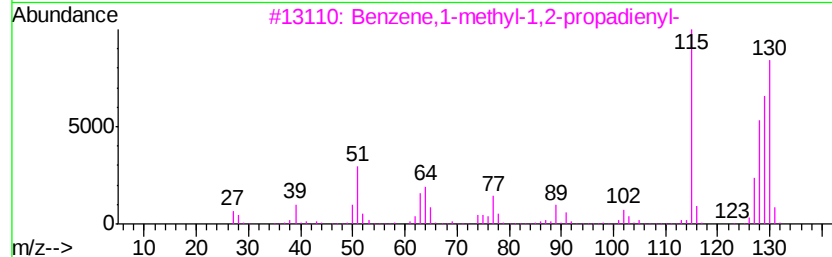
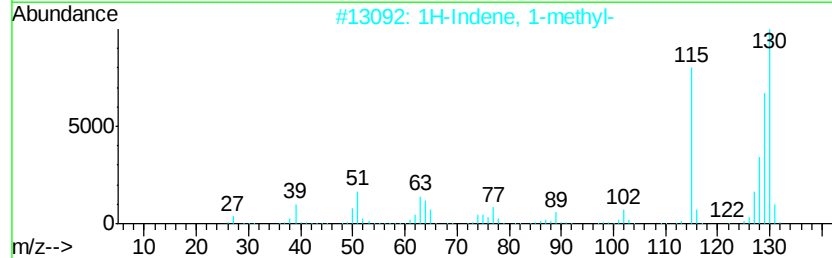
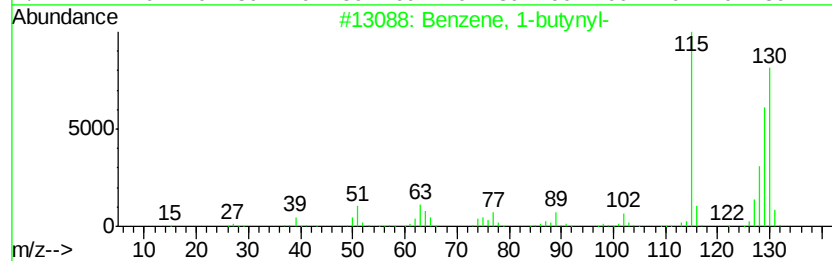
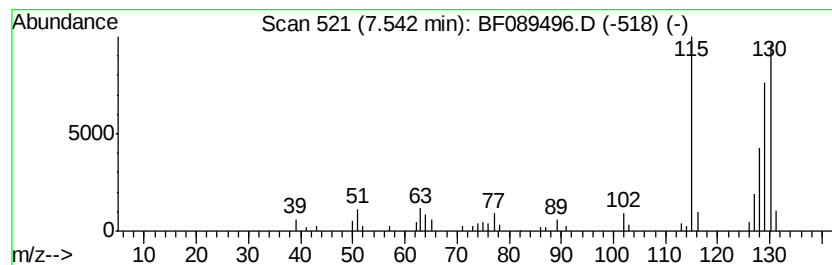
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzene, 1-butynyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.54	4.36 ng	82507	Naphthalene-d8	7.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-butynyl-	130	C10H10	000622-76-4	95
2			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
3			Benzene, 1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
4			2-Methylindene	130	C10H10	002177-47-1	91
5			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF073116\
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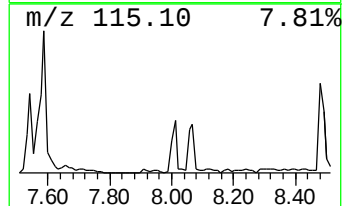
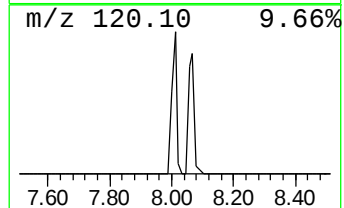
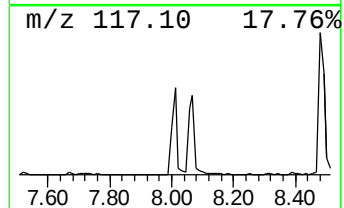
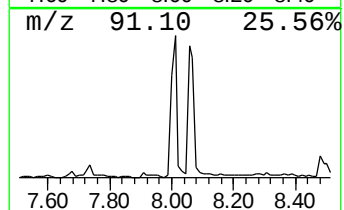
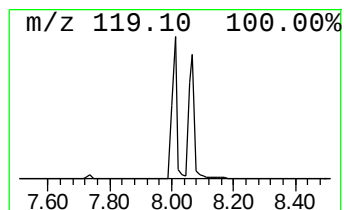
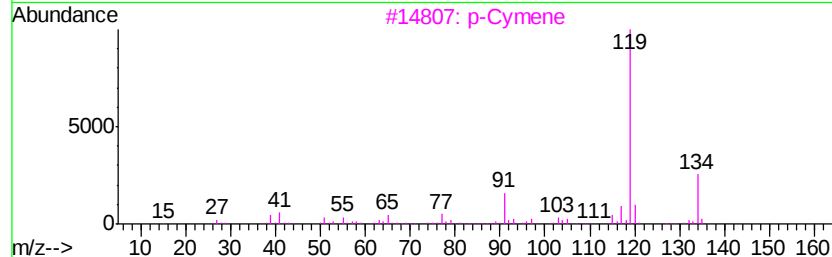
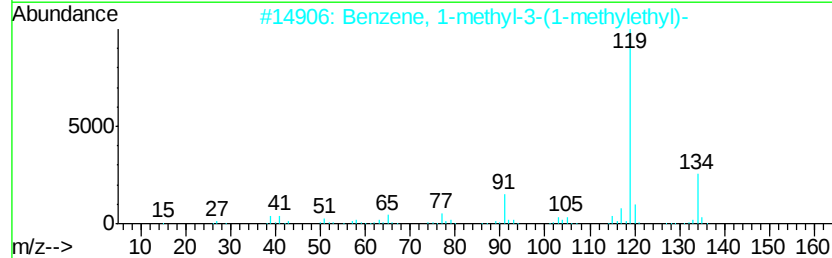
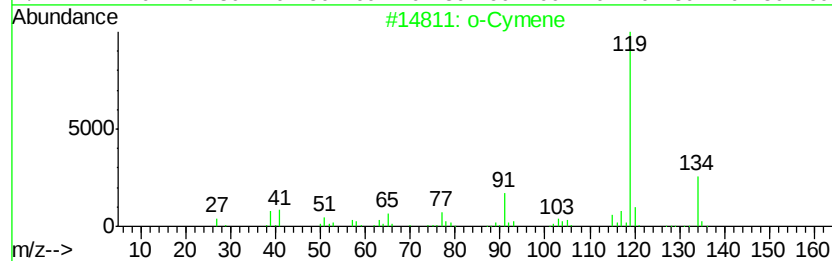
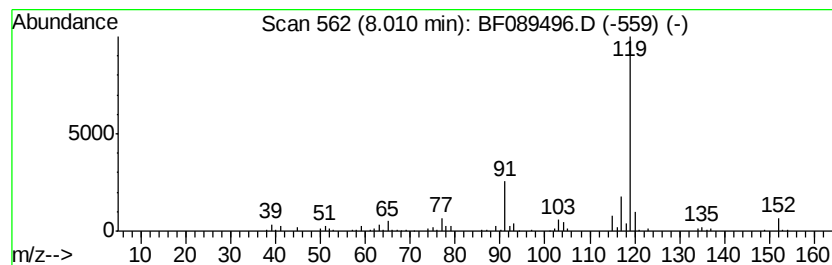
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.01	18.86 ng	356633	Napthalene-d8	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	87
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	87
3		p-Cymene	134	C10H14	000099-87-6	86
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	80
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	80



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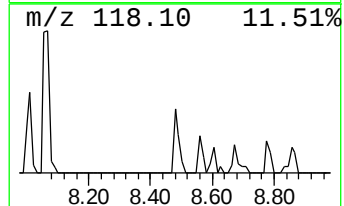
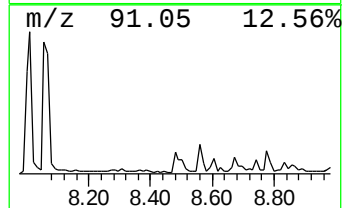
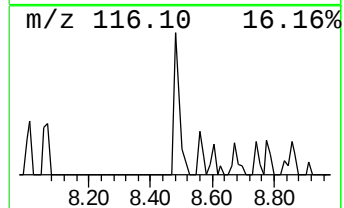
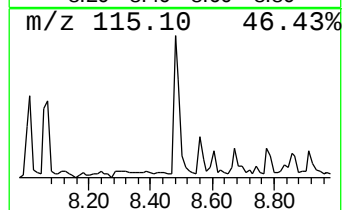
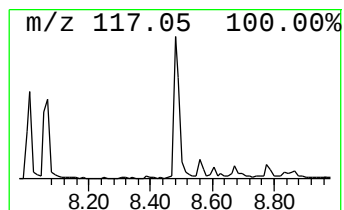
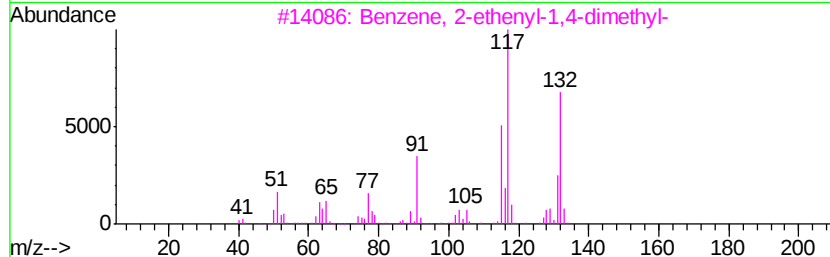
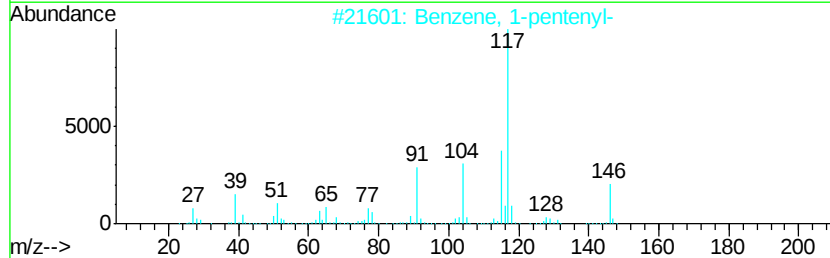
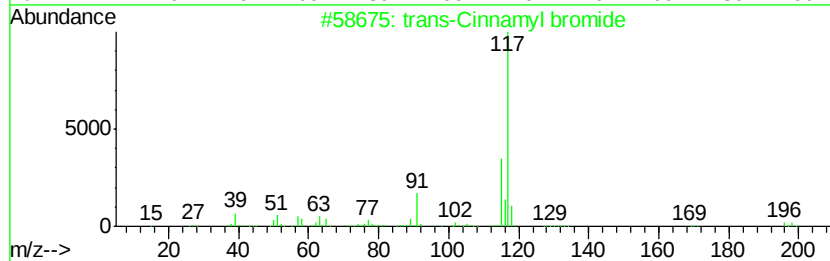
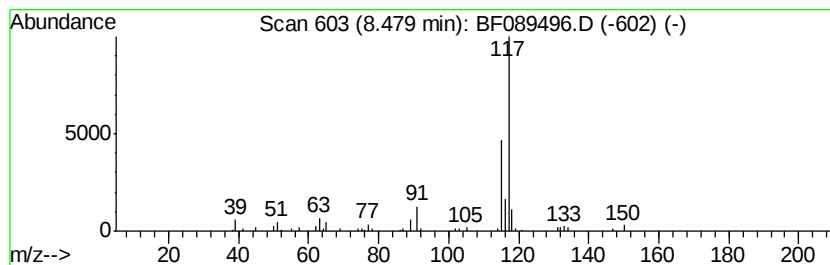
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 trans-Cinnamyl bromide Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.48	7.00 ng	132443	Naphthalene-d8	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	64
2		Benzene, 1-pentenyl-	146	C11H14	000826-18-6	64
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	64
4		Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
5		2,4-Dimethylstyrene	132	C10H12	002234-20-0	56



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
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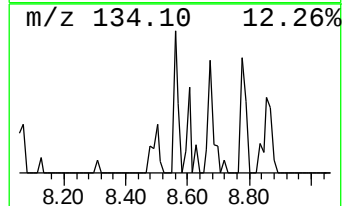
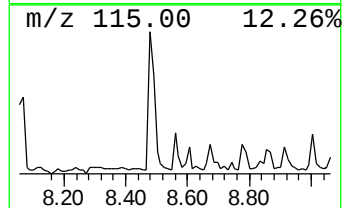
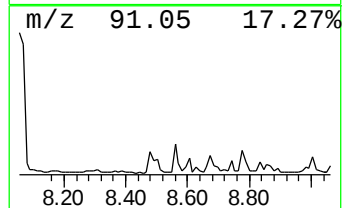
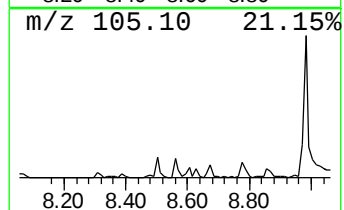
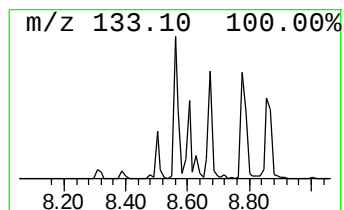
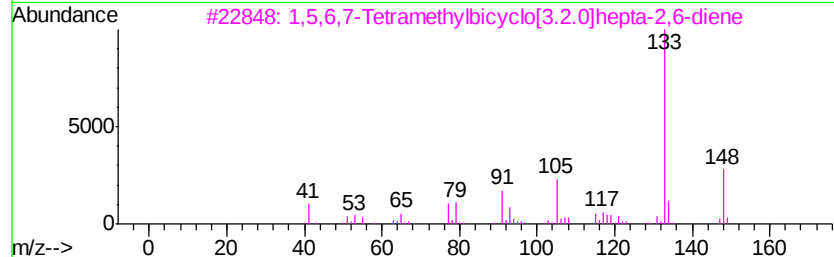
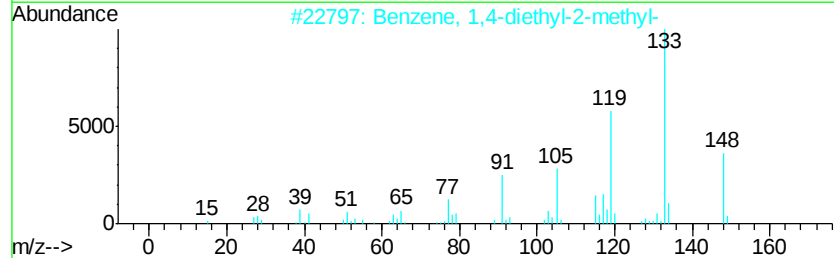
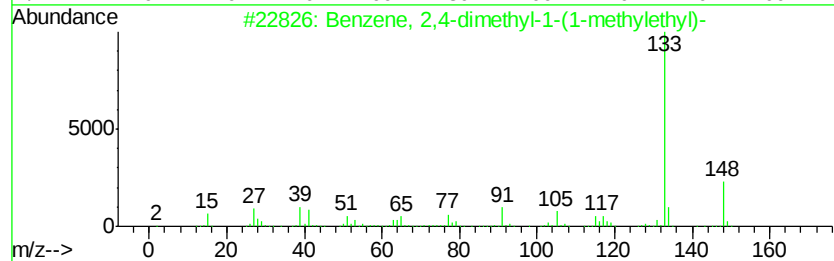
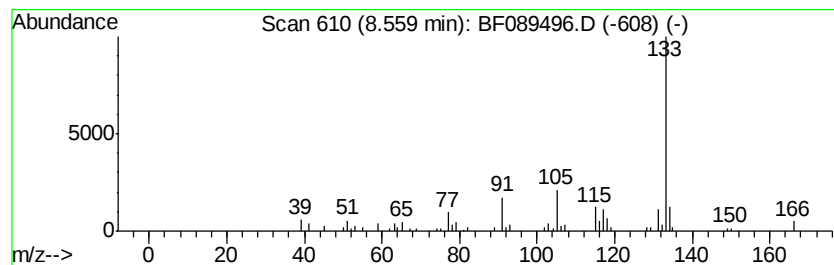
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Benzene, 2,4-dimethyl-1-(1-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.56	4.99 ng	94277	Naphthalene-d8	7.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2,4-dimethyl-1-(1-methyl-...	148	C11H16	004706-89-2	78
2			Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	78
3			1,5,6,7-Tetramethylbicyclo[3.2.0]hepta-2,6-diene	148	C11H16	134329-46-7	64
4			Benzene, 1,3-dimethyl-5-(1-methyl-...	148	C11H16	004706-90-5	64
5			3,4-Dimethylcumene	148	C11H16	1000370-34-1	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
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Misc :
ALS Vial : 31 Sample Multiplier: 1

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ClientSampleId :
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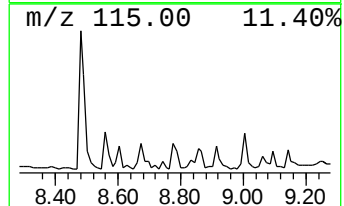
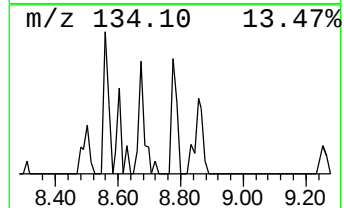
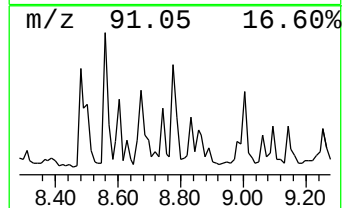
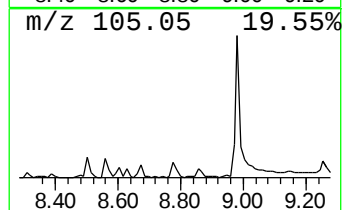
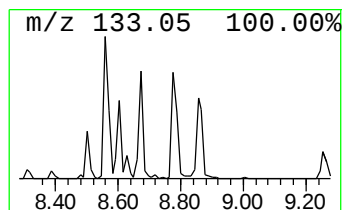
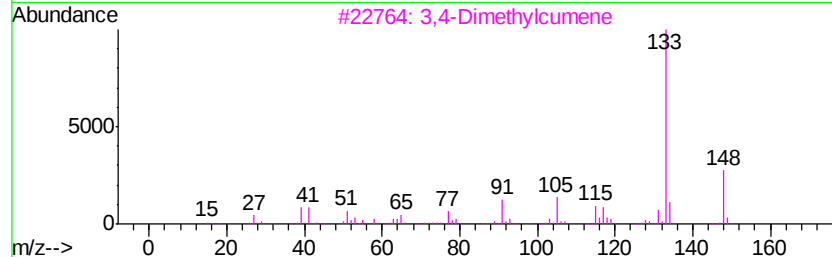
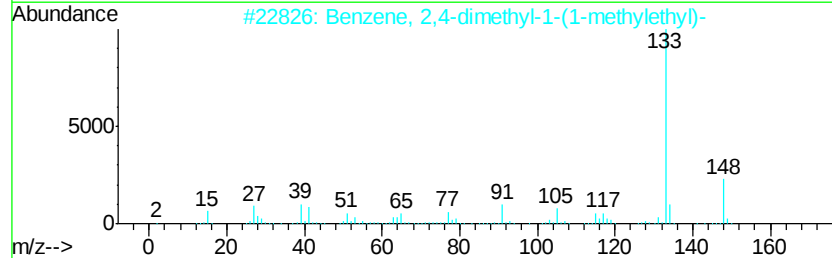
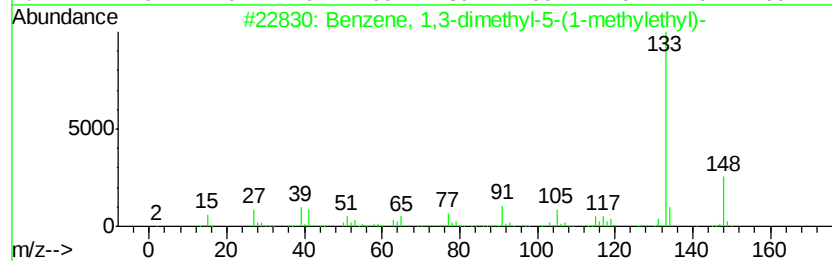
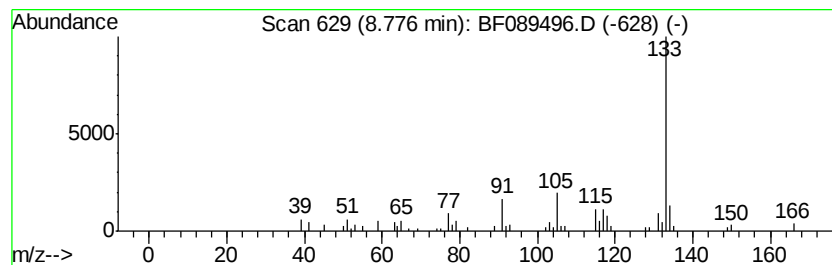
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzene, 1,3-dimethyl-5-(1-... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.78	4.06 ng	78729	Acenaphthene-d10	9.52

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-5-(1-methyl-...	148	C11H16	004706-90-5	78
2			Benzene, 2,4-dimethyl-1-(1-methyl-...	148	C11H16	004706-89-2	78
3			3,4-Dimethylcumene	148	C11H16	1000370-34-1	72
4			Benzene, 1,4-dimethyl-2-(1-methyl-...	148	C11H16	004132-72-3	64
5			Oxalamide, N-(2-hydroxyethyl)-N'...	278	C14H18N2O4	1000304-26-2	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF073116\
Data File : BF089496.D
Acq On : 1 Aug 2016 5:31
Operator : UM/SJ
Sample : H4215-04
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-14-COMP

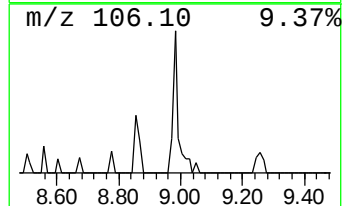
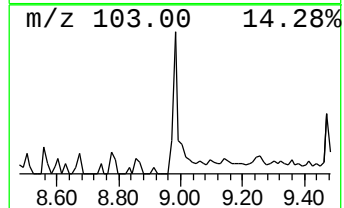
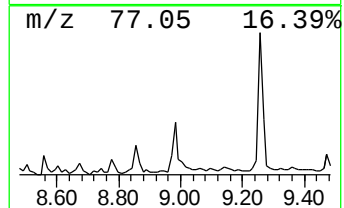
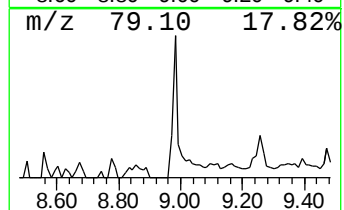
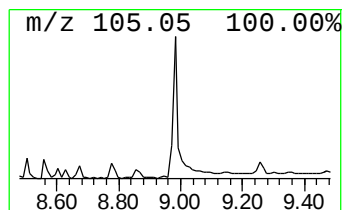
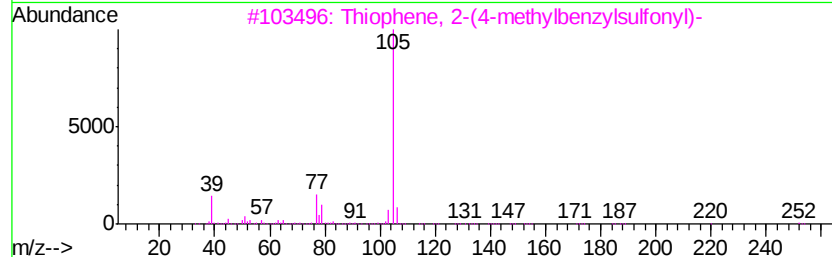
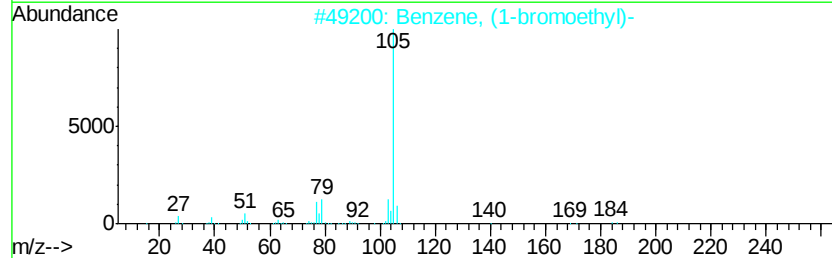
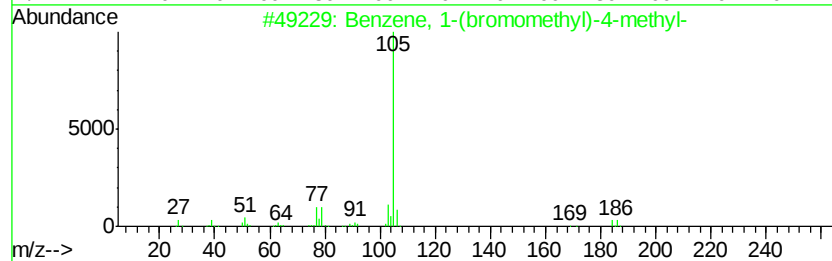
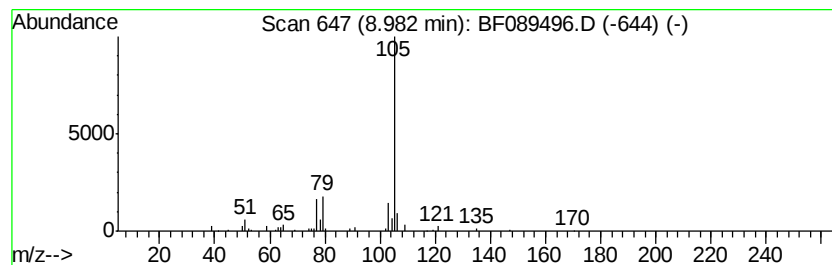
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Benzene, 1-(bromomethyl)-4-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	8.56 ng	166009	Acenaphthene-d10	9.52

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-(bromomethyl)-4-methyl-	184	C8H9Br	000104-81-4	80
2			Benzene, (1-bromoethyl)-	184	C8H9Br	000585-71-7	72
3			Thiophene, 2-(4-methylbenzylsulf...	252	C12H12O2S2	153837-88-8	64
4			Benzene, (1-nitroethyl)-	151	C8H9NO2	007214-61-1	59
5			3-Phenylpropanoic anhydride	282	C18H18O3	1000333-89-6	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF073116\
Data File : BF089496.D
Acq On : 1 Aug 2016 5:31
Operator : UM/SJ
Sample : H4215-04
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-14-COMP

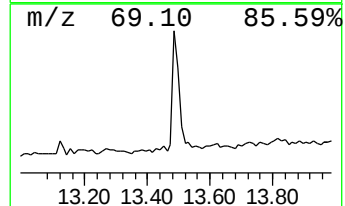
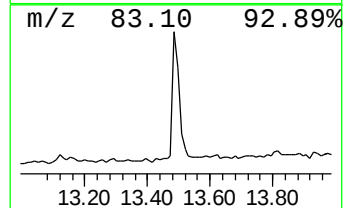
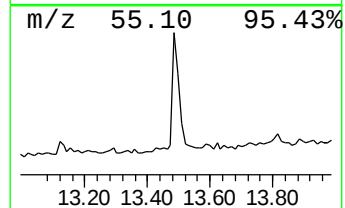
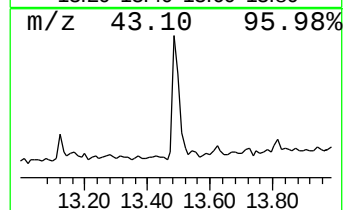
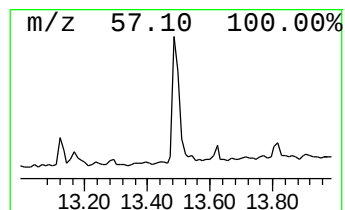
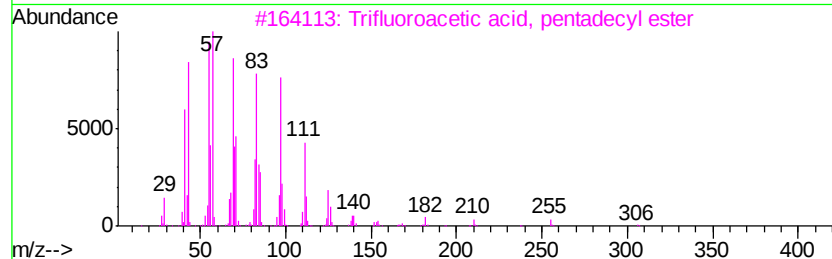
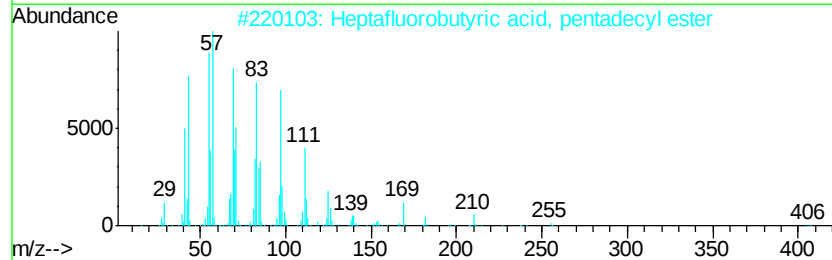
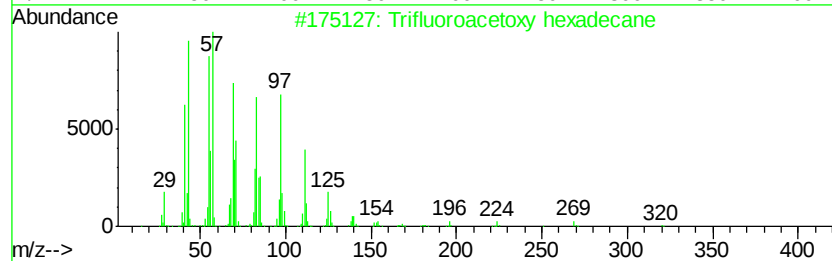
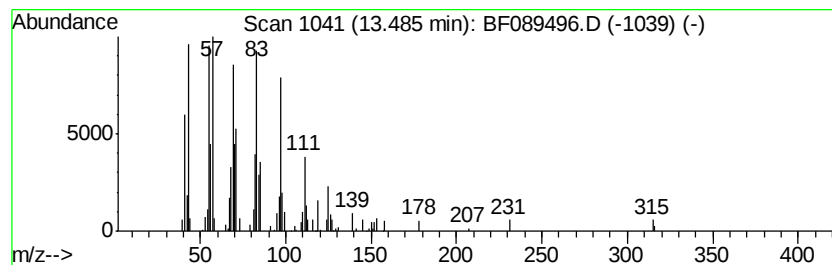
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Trifluoroacetoxy hexadecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	5.73 ng	112756	Chrysene-d12	13.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	95
2		Heptafluorobutyric acid, pentadecyl ester	424	C19H31F7O2	959261-23-5	95
3		Trifluoroacetic acid, pentadecyl ester	324	C17H31F3O2	959010-23-2	95
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	95
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	95



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF073116\
 Data File : BF089496.D
 Acq On : 1 Aug 2016 5:31
 Operator : UM/SJ
 Sample : H4215-04
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-14-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.58	31.4	ng	457198	1	6.49	291475	20.0
Bicyclo[4.2.0]oct...	5.27	11.1	ng	161032	1	6.49	291475	20.0
unknown6.26	6.26	90.8	ng	1323000	1	6.49	291475	20.0
Benzene, 1-etheny...	6.33	19.2	ng	280278	1	6.49	291475	20.0
Indene	6.75	8.5	ng	124145	1	6.49	291475	20.0
Benzene, 2-etheny...	7.13	8.7	ng	127103	1	6.49	291475	20.0
3,4-Dimethylthiop...	7.36	19.9	ng	376588	2	7.78	378233	20.0
Benzene, 1-butynyl-	7.54	4.4	ng	82507	2	7.78	378233	20.0
o-Cymene	8.01	18.9	ng	356633	2	7.78	378233	20.0
trans-Cinnamyl br...	8.48	7.0	ng	132443	2	7.78	378233	20.0
Benzene, 2,4-dime...	8.56	5.0	ng	94277	2	7.78	378233	20.0
Benzene, 1,3-dime...	8.78	4.1	ng	78729	3	9.52	387713	20.0
Benzene, 1-(bromo...	8.98	8.6	ng	166009	3	9.52	387713	20.0
Trifluoroacetoxy ...	13.49	5.7	ng	112756	5	13.61	393598	20.0