

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
 Data File : BF085182.D
 Acq On : 4 Mar 2016 3:17
 Operator : SJ/UM
 Sample : H1648-08
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-49(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-BF030316.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	5.190	251	254	260	rVB	688145	940788	22.91%	2.231%
2	5.590	286	289	291	rBV	3811223	4106251	100.00%	9.735%
3	6.596	374	377	380	rBV	2769914	4006237	97.56%	9.498%
4	6.744	387	390	393	rBV	3959601	3836058	93.42%	9.095%
5	6.973	408	410	413	rBV	830293	895510	21.81%	2.123%
6	7.133	422	424	427	rVB	3207026	3017249	73.48%	7.154%
7	7.544	456	460	462	rBV	2083831	2853805	69.50%	6.766%
8	7.613	462	466	468	rVB	58870	65961	1.61%	0.156%
9	8.264	520	523	525	rBV	1363998	1295802	31.56%	3.072%
10	9.339	614	617	619	rBV	4417014	3972503	96.74%	9.418%
11	9.727	649	651	653	rBV	741521	601000	14.64%	1.425%
12	9.762	653	654	656	rVB	59169	63168	1.54%	0.150%
13	9.945	668	670	672	rVB	137529	109255	2.66%	0.259%
14	10.013	674	676	678	rVB2	1061497	1270329	30.94%	3.012%
15	10.173	687	690	692	rBV	25366	44054	1.07%	0.104%
16	10.448	710	714	715	rBV	223073	238553	5.81%	0.566%
17	10.562	722	724	727	rVB	46653	47798	1.16%	0.113%
18	10.665	730	733	736	rBV	71956	119337	2.91%	0.283%
19	10.745	739	740	742	rBV	29736	49382	1.20%	0.117%
20	10.802	742	745	748	rVB	2357183	2428477	59.14%	5.758%
21	10.916	753	755	762	rVB	257171	313429	7.63%	0.743%
22	11.110	770	772	774	rBV	44842	59741	1.45%	0.142%
23	11.362	792	794	796	rBV	126166	119739	2.92%	0.284%
24	11.396	796	797	800	rVB2	56808	50127	1.22%	0.119%
25	11.499	804	806	810	rBV2	1377689	1622740	39.52%	3.847%
26	11.579	810	813	814	rVB2	64915	103148	2.51%	0.245%
27	12.025	848	852	858	rBV3	436183	561519	13.67%	1.331%
28	12.116	858	860	864	rVB2	70078	103417	2.52%	0.245%
29	12.299	873	876	880	rBV3	28633	65979	1.61%	0.156%
30	12.619	902	904	908	rVB	31786	54046	1.32%	0.128%
31	12.711	910	912	915	rBV	501955	442044	10.77%	1.048%
32	12.802	919	920	924	rVV3	77568	120595	2.94%	0.286%
33	12.939	928	932	935	rBV	410574	459363	11.19%	1.089%
34	13.088	942	945	947	rBV	3464115	3246236	79.06%	7.696%

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35	13.271	959	961	964	rVB	40883	52630	1.28%	0.125%
36	13.979	1021	1023	1027	rVB	240276	331258	8.07%	0.785%
37	14.139	1033	1037	1038	rBV2	904195	1197308	29.16%	2.839%
38	14.162	1038	1039	1041	rVB	137910	105343	2.57%	0.250%
39	14.608	1075	1078	1080	rBV3	63634	100956	2.46%	0.239%
40	14.894	1101	1103	1106	rBV2	85709	147924	3.60%	0.351%
41	14.985	1109	1111	1116	rVB	90838	116801	2.84%	0.277%
42	15.179	1125	1128	1132	rBV	104943	246496	6.00%	0.584%
43	15.248	1132	1134	1136	rBV	71053	91042	2.22%	0.216%
44	15.477	1152	1154	1157	rVB	76782	105481	2.57%	0.250%
45	15.545	1157	1160	1162	rBV	79540	117482	2.86%	0.279%
46	15.602	1162	1165	1167	rBV	546218	788829	19.21%	1.870%
47	16.071	1204	1206	1209	rBV	51106	83346	2.03%	0.198%
48	16.128	1209	1211	1219	rVB8	33774	112257	2.73%	0.266%
49	16.322	1225	1228	1234	rVB3	87513	197133	4.80%	0.467%
50	16.517	1240	1245	1248	rBV7	29329	94762	2.31%	0.225%
51	16.585	1248	1251	1254	rBV	47469	96530	2.35%	0.229%
52	16.780	1264	1268	1271	rVB6	22875	58325	1.42%	0.138%
53	17.065	1290	1293	1296	rBV2	46862	93228	2.27%	0.221%
54	17.134	1296	1299	1302	rBV2	73319	137875	3.36%	0.327%
55	17.271	1307	1311	1317	rBV5	25196	92160	2.24%	0.219%
56	17.442	1322	1326	1329	rBV3	45243	110315	2.69%	0.262%
57	17.511	1329	1332	1338	rVB	42184	93282	2.27%	0.221%
58	17.683	1343	1347	1354	rBV10	41091	146522	3.57%	0.347%
59	17.900	1362	1366	1372	rVB3	32678	104825	2.55%	0.249%
60	18.551	1421	1423	1430	rVB6	27395	72432	1.76%	0.172%

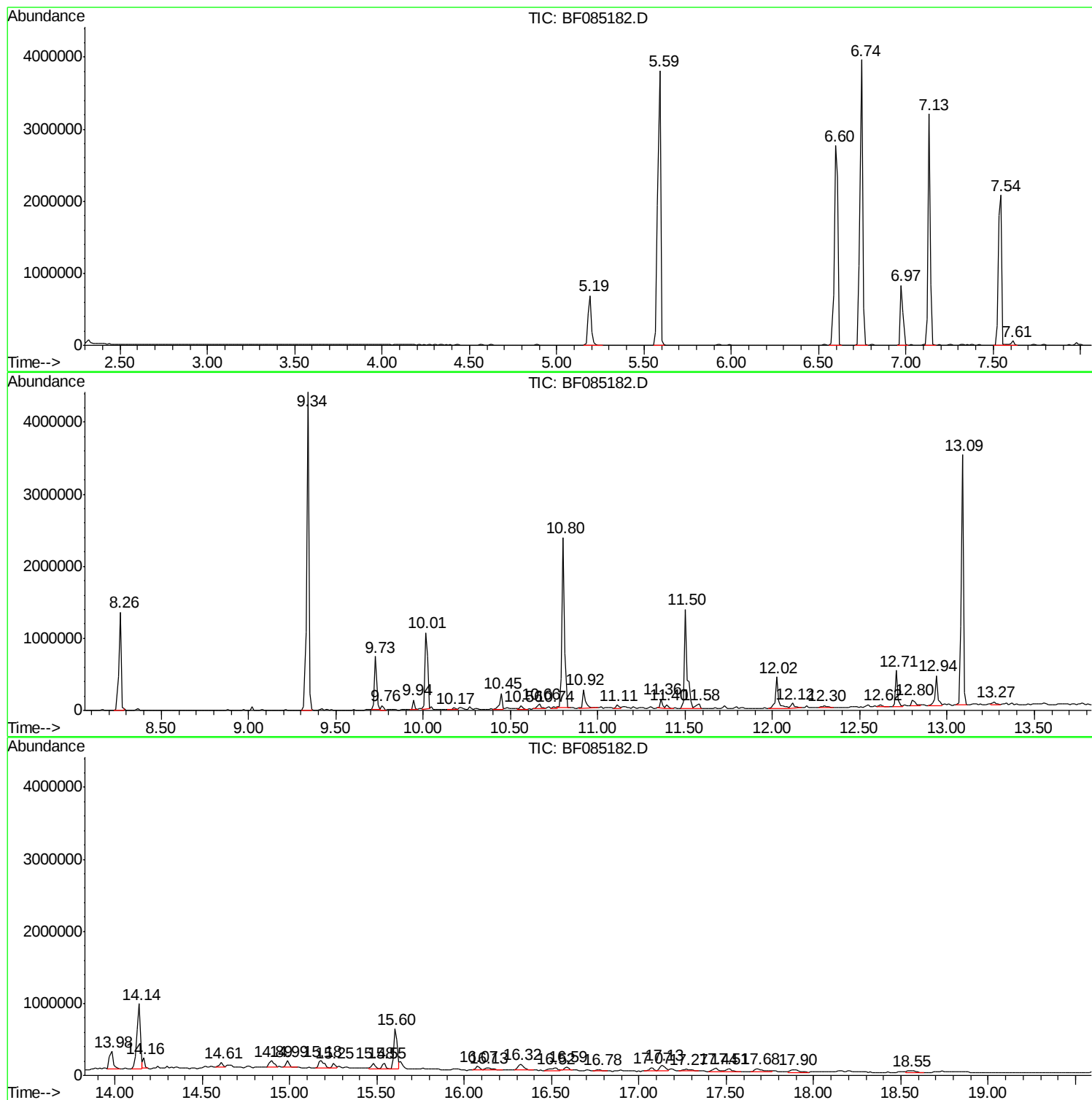
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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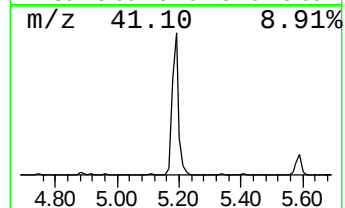
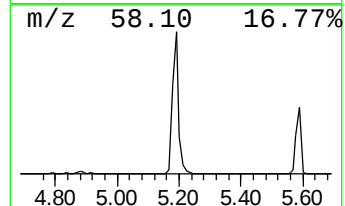
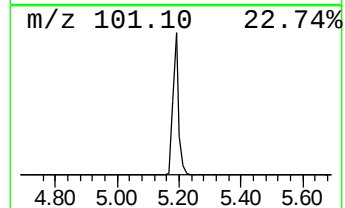
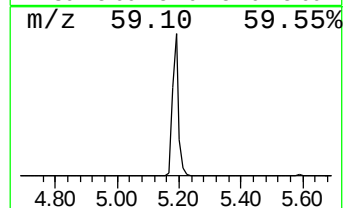
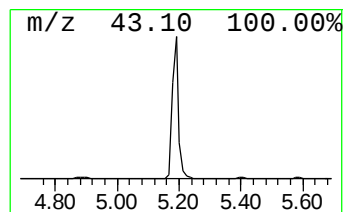
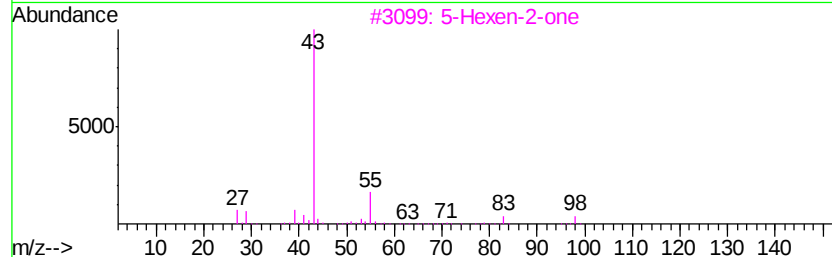
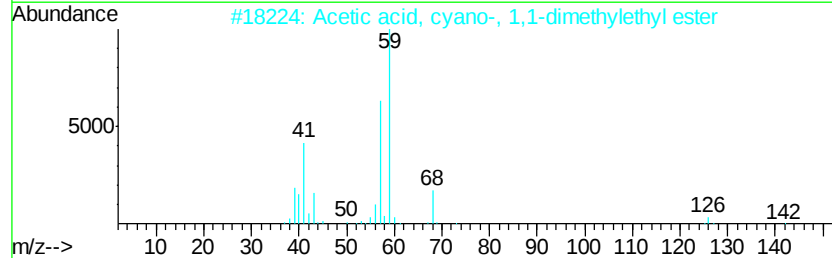
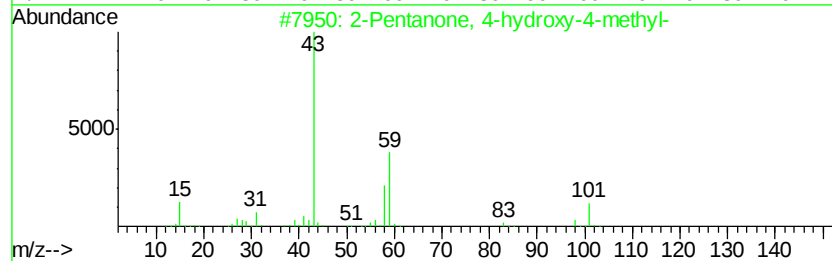
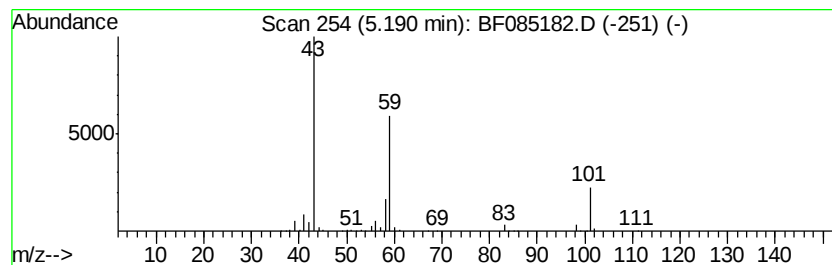
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	21.01 ng	940788	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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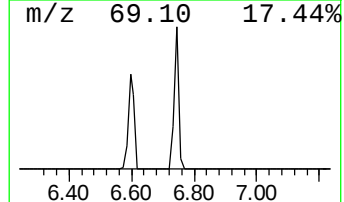
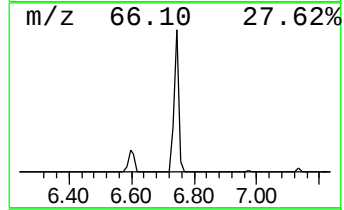
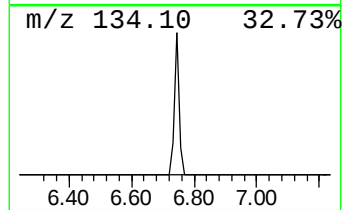
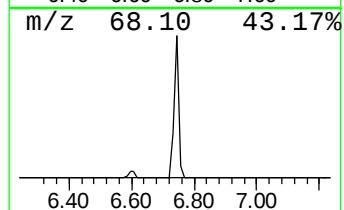
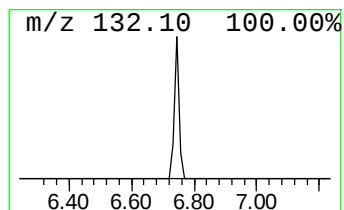
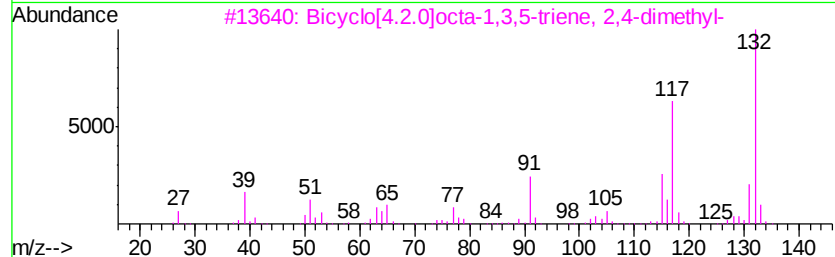
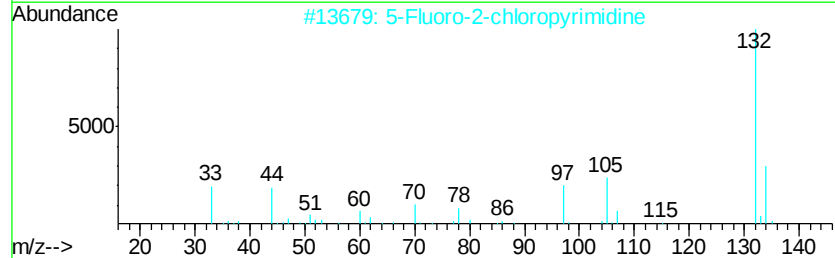
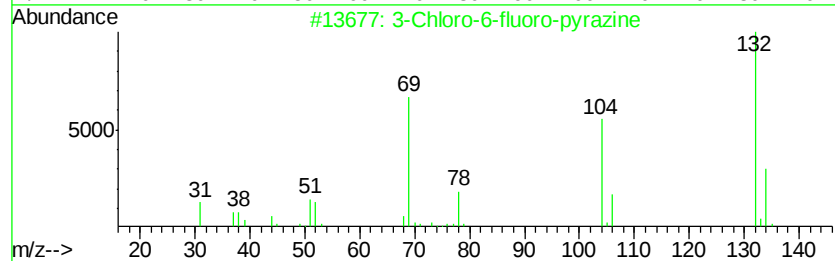
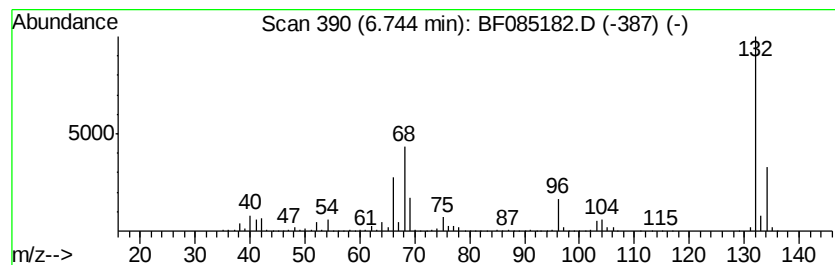
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	85.67 ng	3836060	1,4-Dichlorobenzene-d4	6.97

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		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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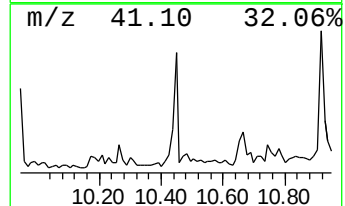
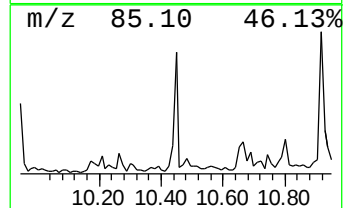
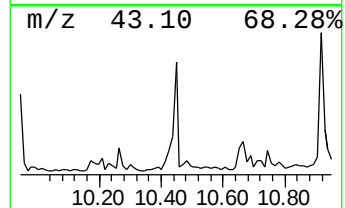
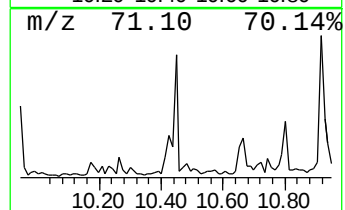
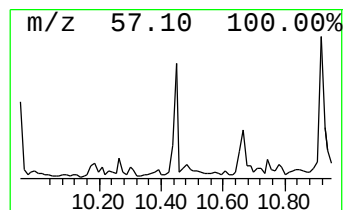
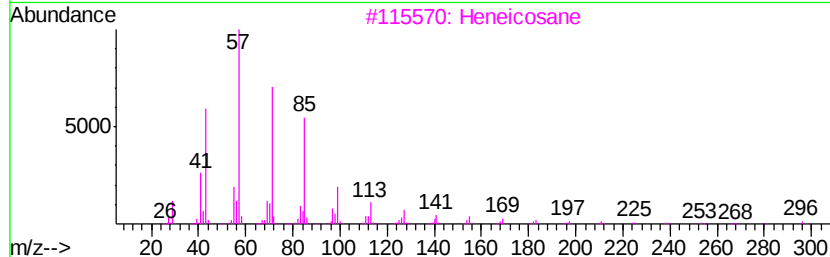
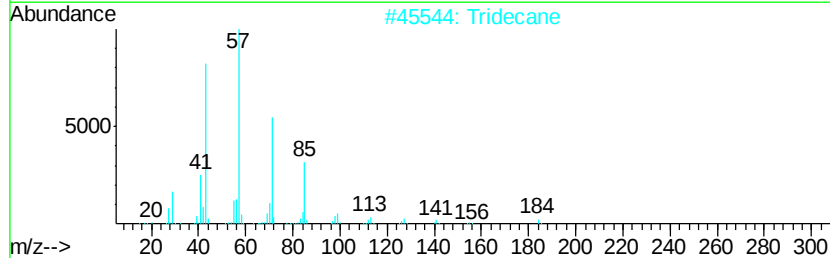
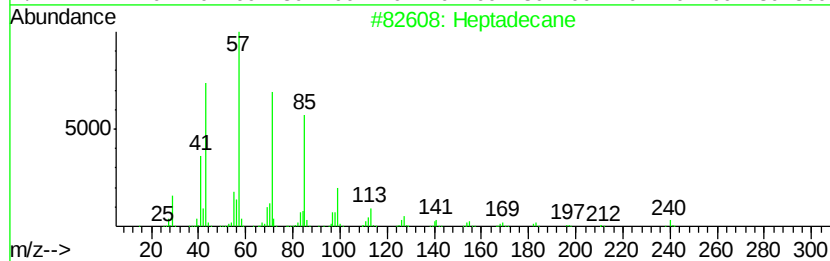
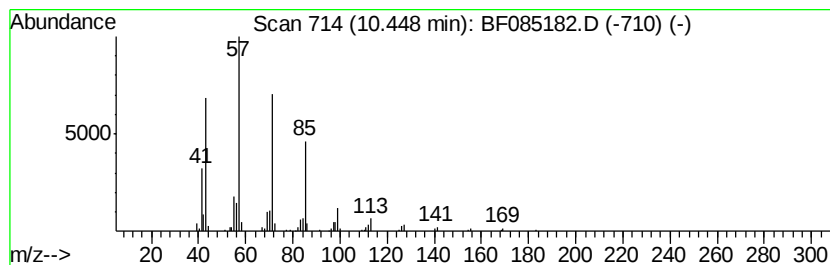
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Heptadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.45	3.76 ng	238553	Acenaphthene-d10	10.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptadecane	240 C17H36	000629-78-7	91
2	Tridecane	184 C13H28	000629-50-5	90
3	Heneicosane	296 C21H44	000629-94-7	90
4	Tetradecane	198 C14H30	000629-59-4	87
5	Dodecane, 1-iodo-	296 C12H25I	004292-19-7	86



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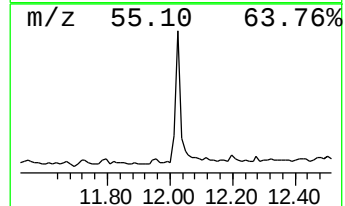
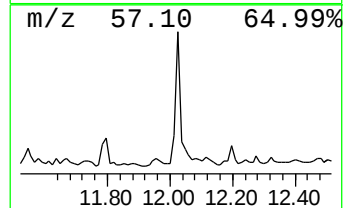
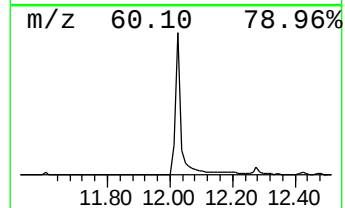
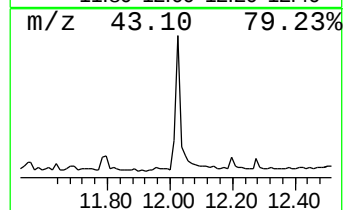
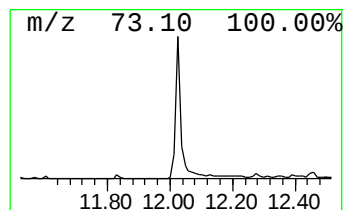
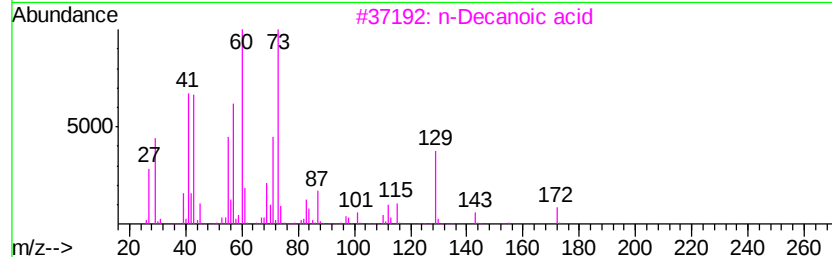
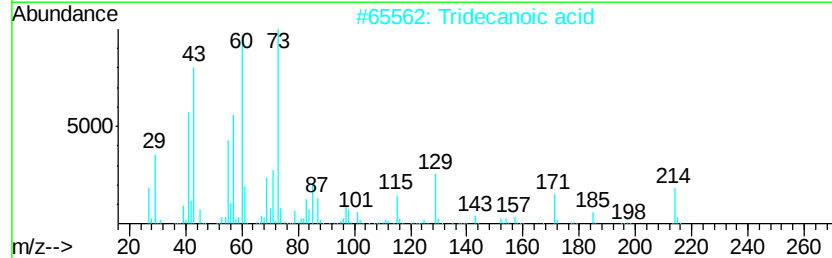
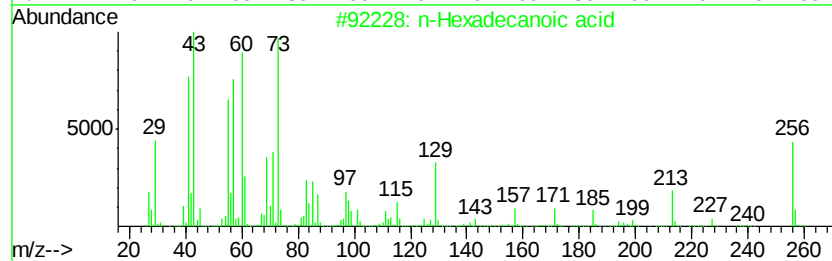
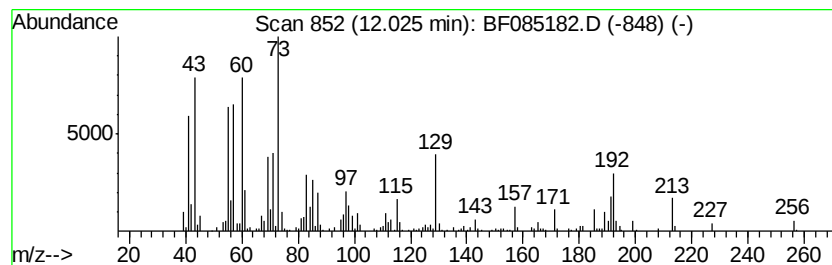
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	6.92 ng	561519	Phenanthrene-d10	11.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tridecanoic acid	214	C13H26O2	000638-53-9	81
3		n-Decanoic acid	172	C10H20O2	000334-48-5	76
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	72
5		Estra-1,3,5(10)-trien-17.beta.-ol	256	C18H24O	002529-64-8	20



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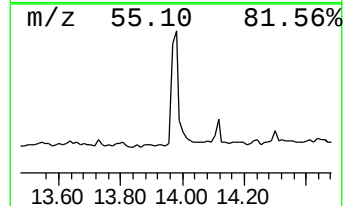
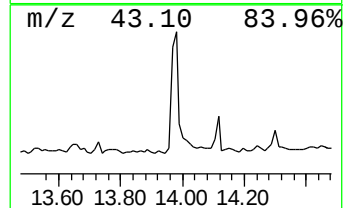
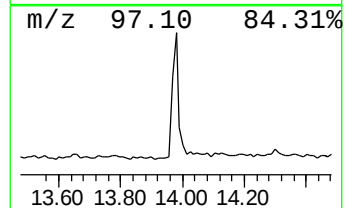
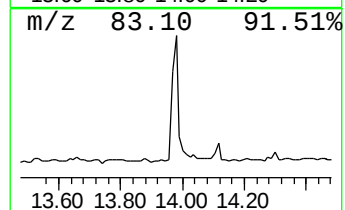
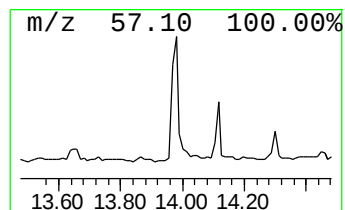
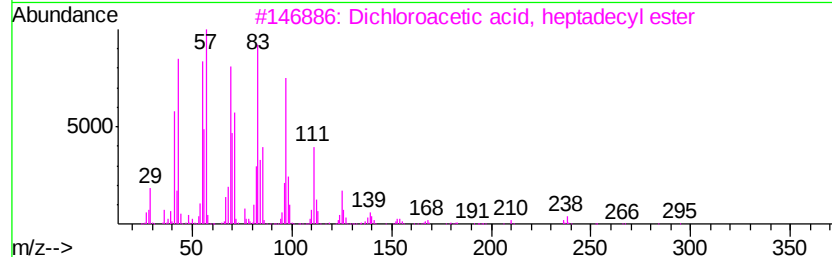
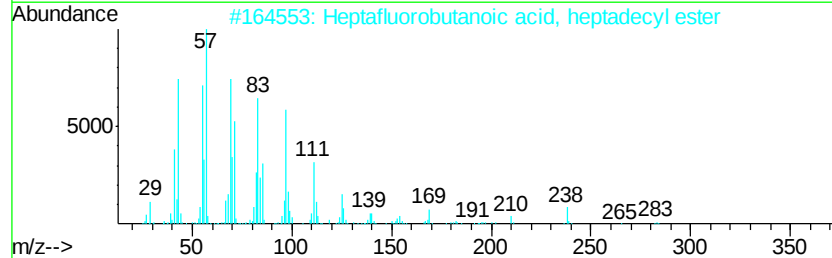
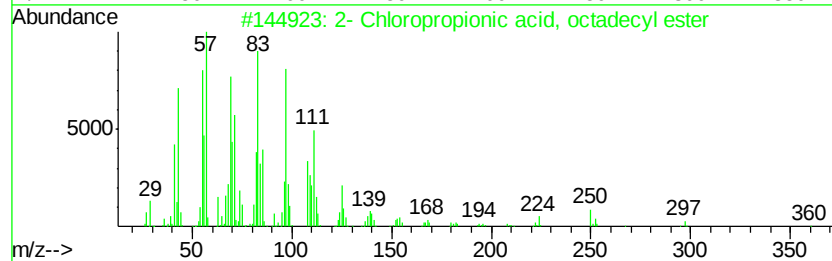
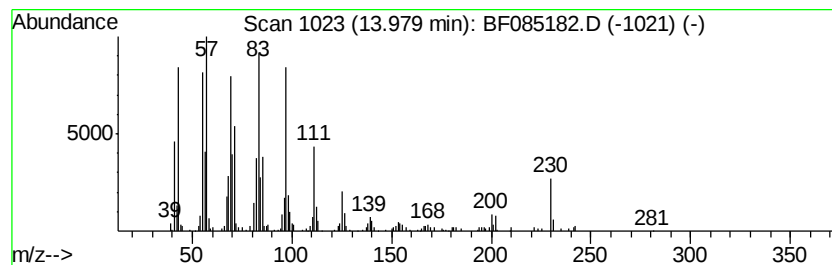
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 2- Chloropropionic acid, oc... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	5.53 ng	331258	Chrysene-d12	14.14

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-	Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
2		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
3		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
4		Cyclopentadecane	210	C15H30	000295-48-7	93
5		1-Nonadecanol	284	C19H40O	001454-84-8	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
Data File : BF085182.D
Acq On : 4 Mar 2016 3:17
Operator : SJ/UM
Sample : H1648-08
Misc :
ALS Vial : 26 Sample Multiplier: 1

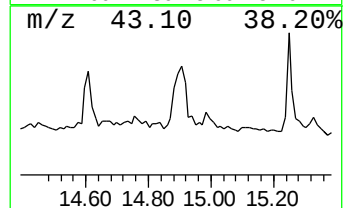
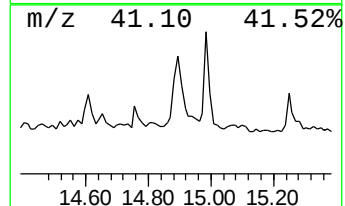
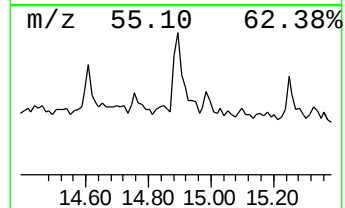
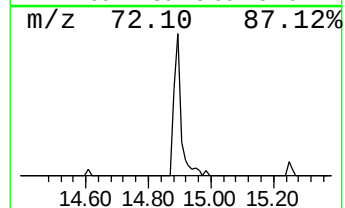
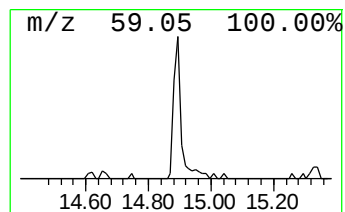
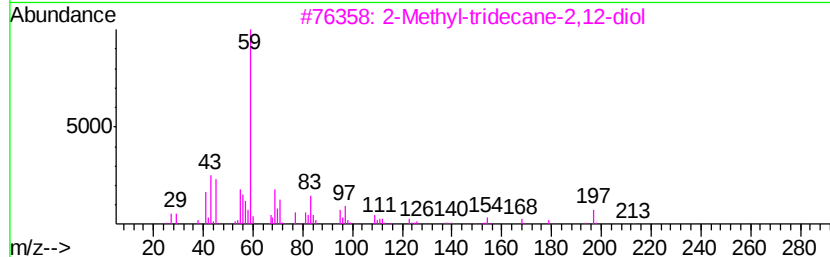
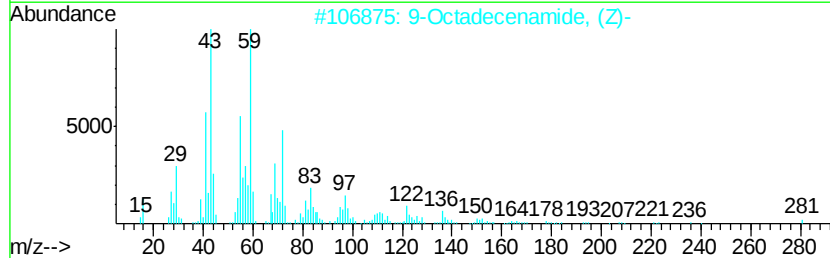
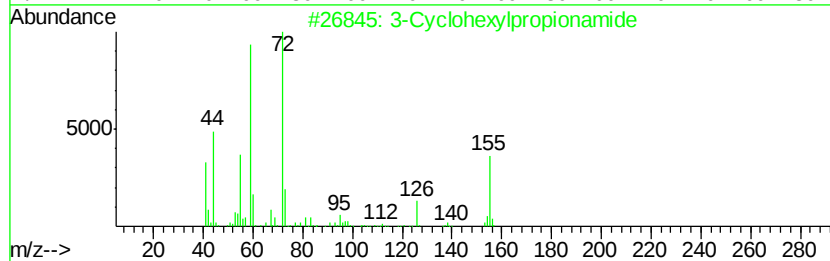
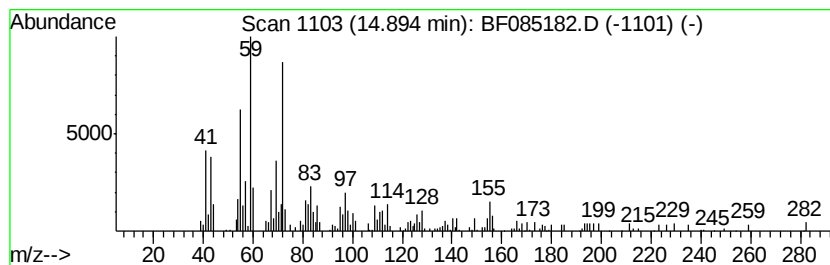
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 3-Cyclohexylpropionamide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.89	3.75 ng	147924	Perylene-d12	15.60	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Cyclohexylpropionamide	155	C9H17NO	004361-29-9	53
2	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	53
3	2-Methyl-tridecane-2,12-diol	230	C14H30O2	1000187-03-5	38
4	3-Tridecanol	200	C13H28O	010289-68-6	35
5	Tetradecanamide	227	C14H29NO	000638-58-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
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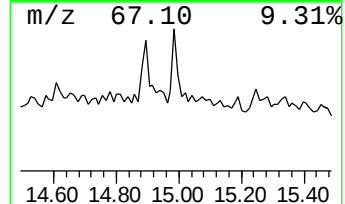
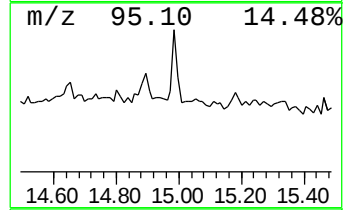
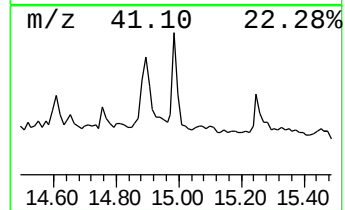
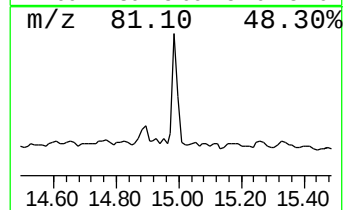
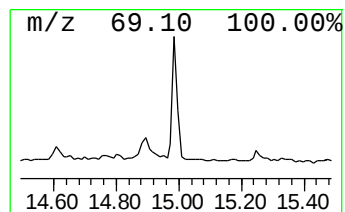
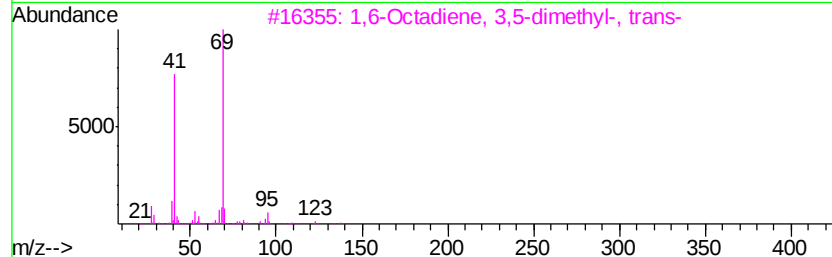
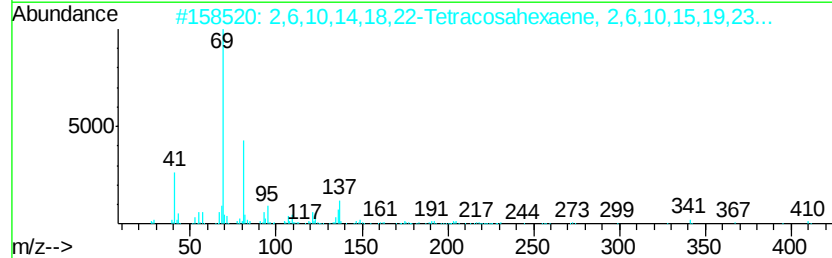
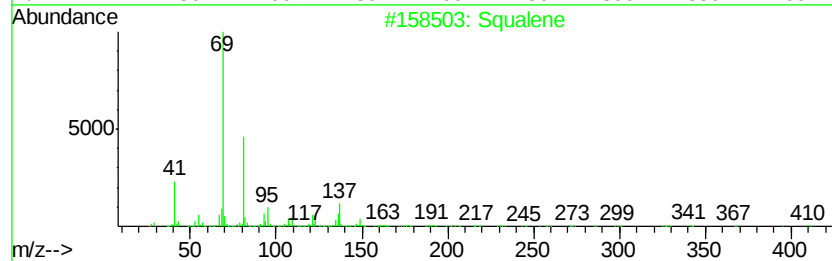
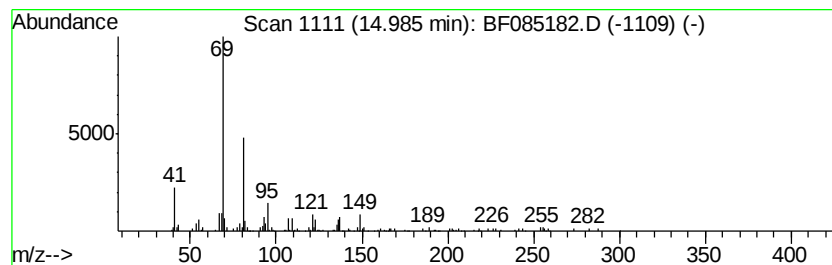
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Squalene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	2.96 ng	116801	Perylene-d12	15.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	72
2		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	64
3		1,6-Octadiene, 3,5-dimethyl-, tr...	138	C10H18	074630-87-8	59
4		4-Oxo-.beta.-damascone	206	C13H18O2	053398-08-6	59
5		2,6-Octadiene, 2,6-dimethyl-	138	C10H18	002792-39-4	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
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Instrument :
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 ClientSampleId :
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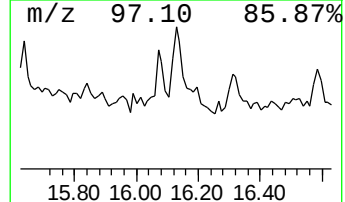
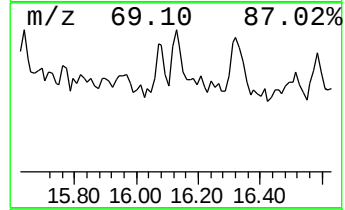
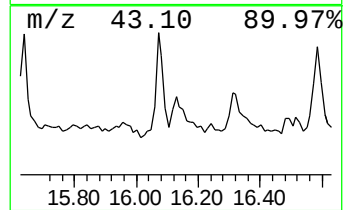
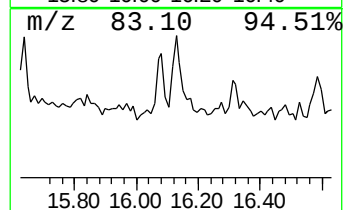
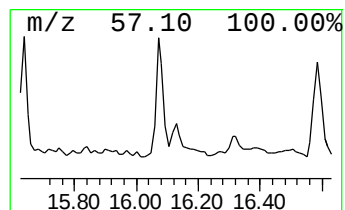
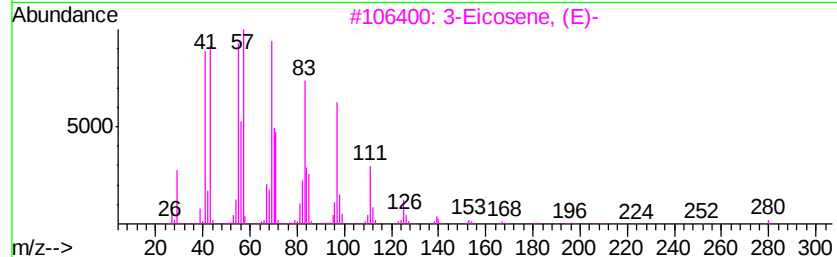
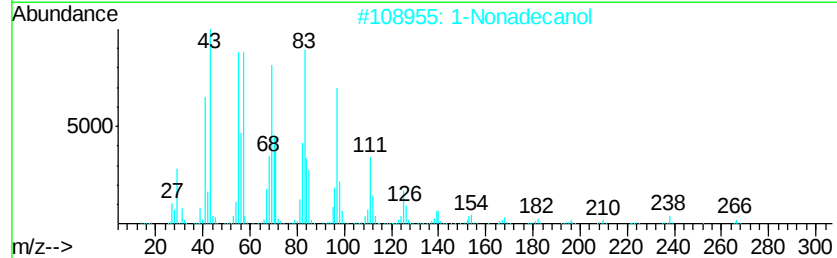
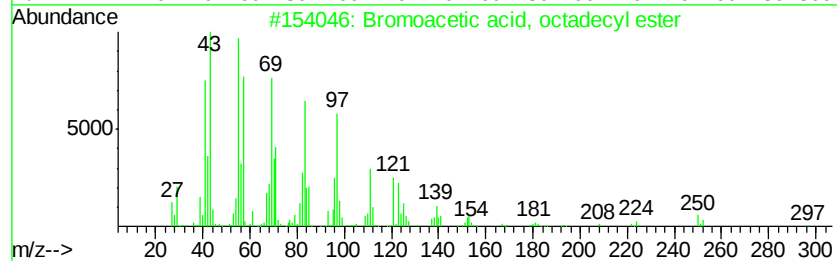
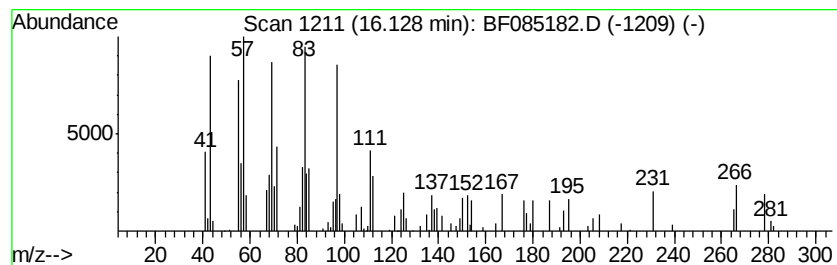
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 11 Bromoacetic acid, octadecyl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	2.85 ng	112257	Perylene-d12	15.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	76
2		1-Nonadecanol	284	C19H40O	001454-84-8	74
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	74
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	74
5		Cyclohexadecane	224	C16H32	000295-65-8	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
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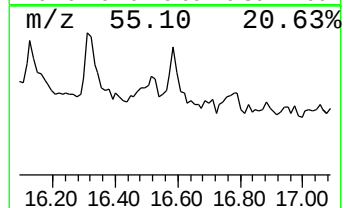
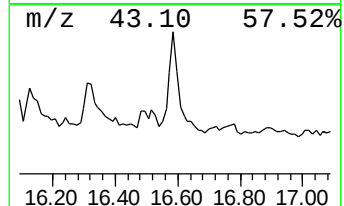
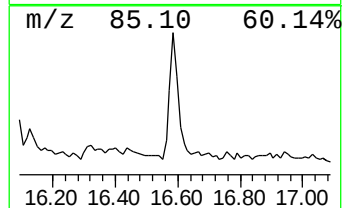
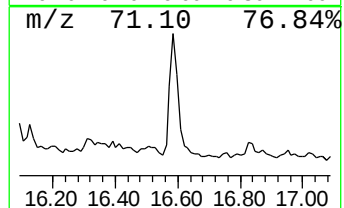
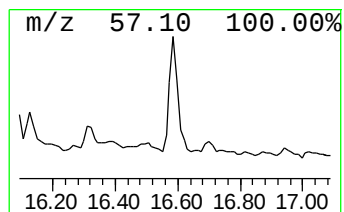
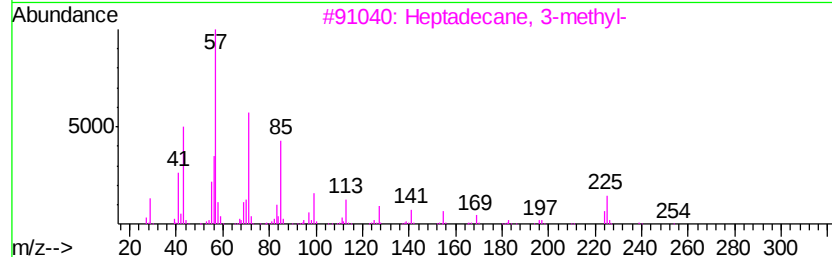
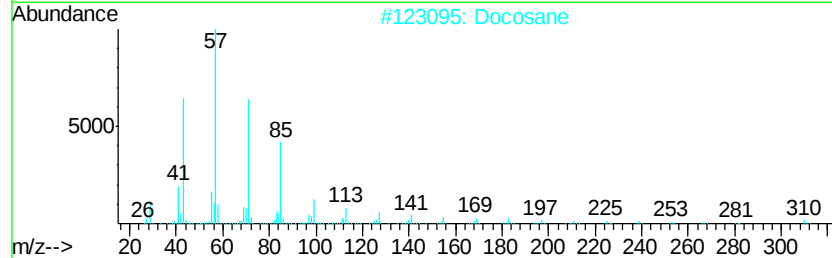
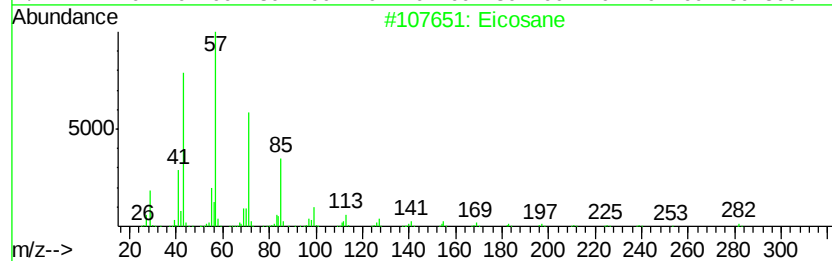
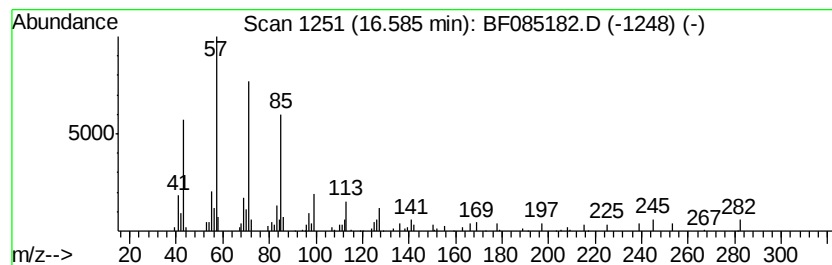
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Eicosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.59	2.45 ng	96530	Perylene-d12	15.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	94
2			Docosane	310	C22H46	000629-97-0	86
3			Heptadecane, 3-methyl-	254	C18H38	006418-44-6	86
4			Heneicosane	296	C21H44	000629-94-7	86
5			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
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Operator : SJ/UM
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Misc :
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ClientSampleId :
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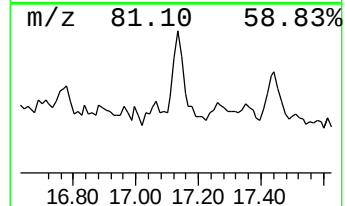
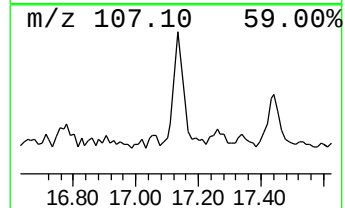
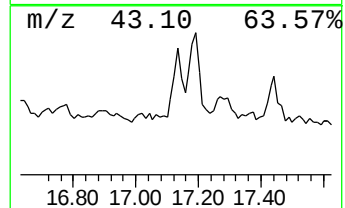
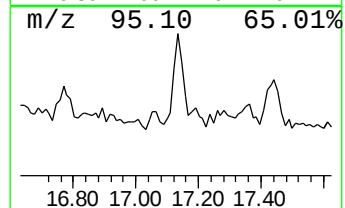
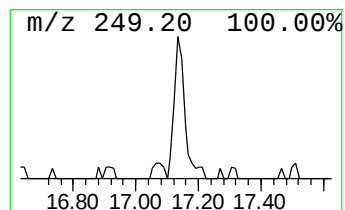
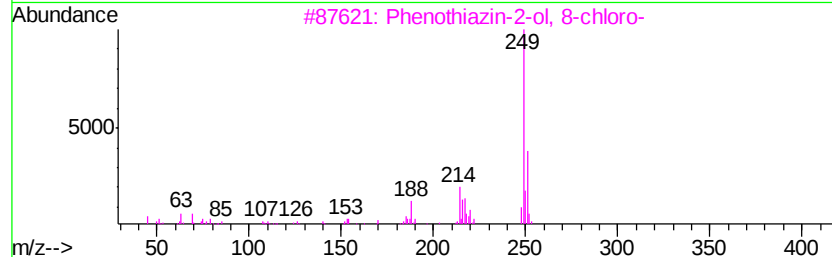
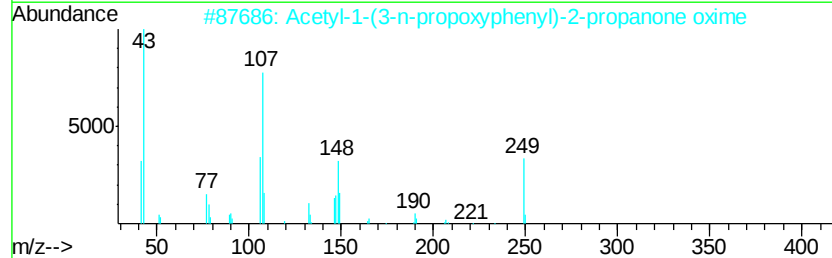
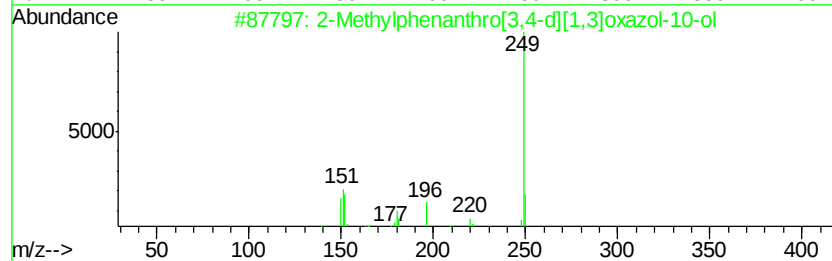
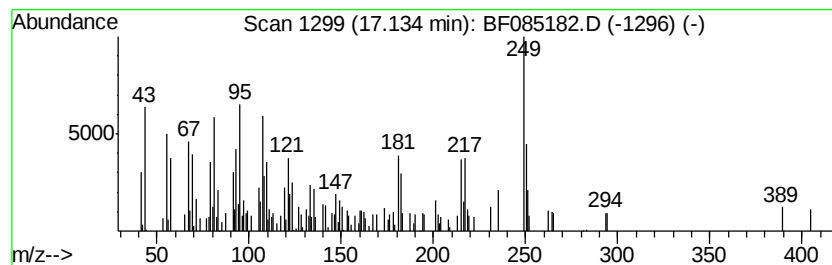
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown17.13 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.13	3.50 ng	137875	Perylene-d12	15.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methylphenanthro[3,4-d][1,3]ox...	249	C16H11N02	098033-24-0	27
2			Acetyl-1-(3-n-propoxyphenyl)-2-p...	249	C14H19N03	1000122-90-2	22
3			Phenothiazin-2-ol, 8-chloro-	249	C12H8ClN0S	005909-61-5	18
4			Isothiocyanic acid, s-phenenyl e...	249	C9H3N3S3	101670-67-1	14
5			7H-Benzo(c)phenothiazine	249	C16H11NS	000226-06-2	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
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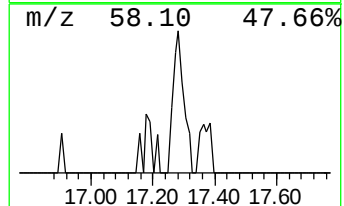
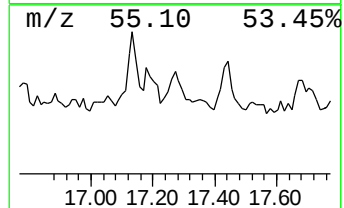
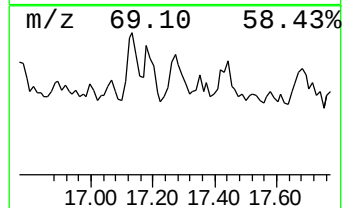
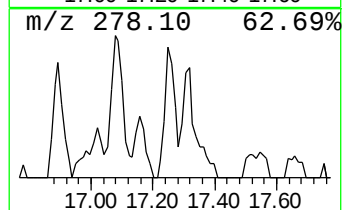
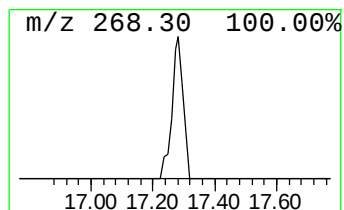
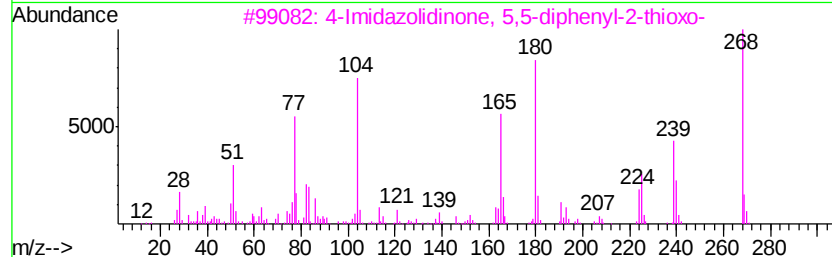
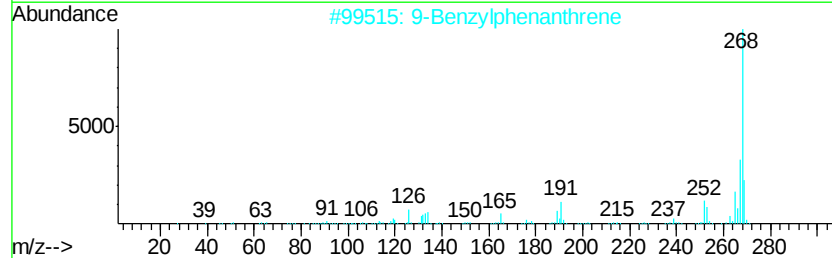
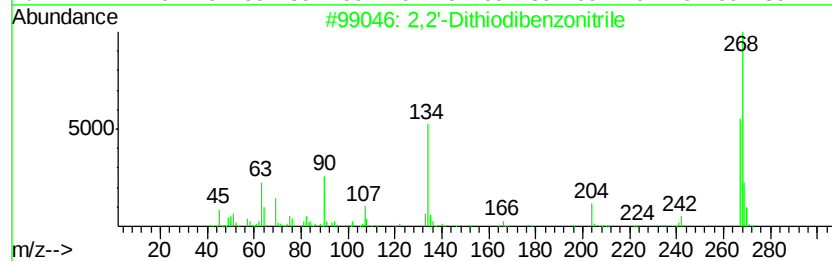
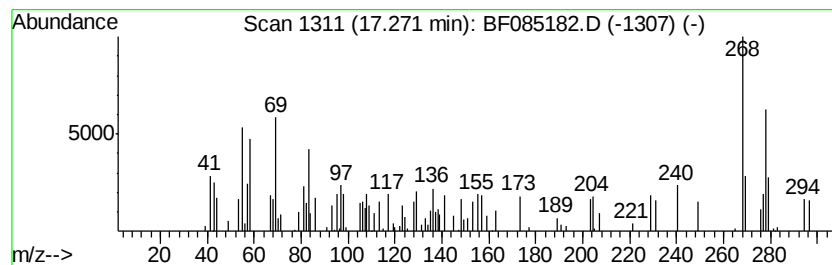
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown17.27 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.27	2.34 ng	92160	Perylene-d12	15.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2'-Dithiodibenzonitrile			268	C14H8N2S2	033174-74-2	30
2	9-Benzylphenanthrene			268	C21H16	000605-05-0	27
3	4-Imidazolidinone, 5,5-diphenyl-...			268	C15H12N2OS	021083-47-6	27
4	1-Thioflavone, 7-methoxy-			268	C16H12O2S	000789-78-6	25
5	4,6-Dibromobenzene-1,3-diol			266	C6H4Br2O2	061524-51-4	22



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Data File : BF085182.D
Acq On : 4 Mar 2016 3:17
Operator : SJ/UM
Sample : H1648-08
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-49(COMP)

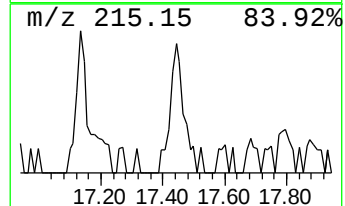
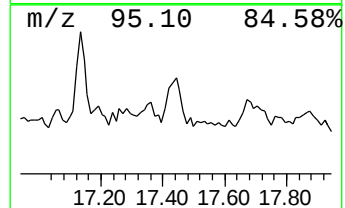
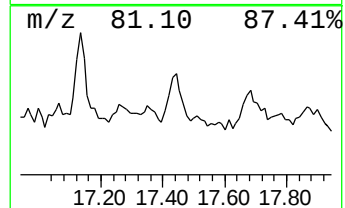
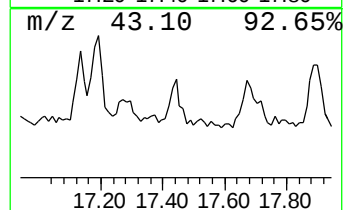
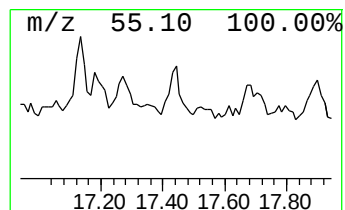
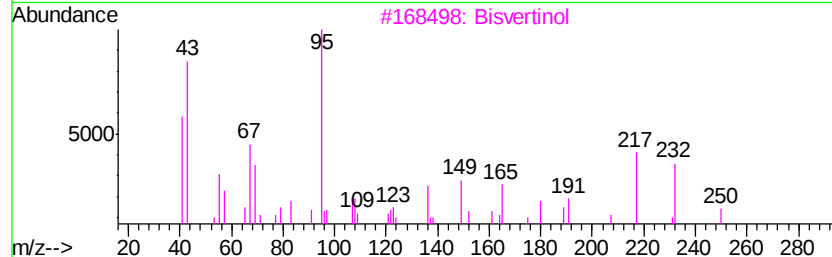
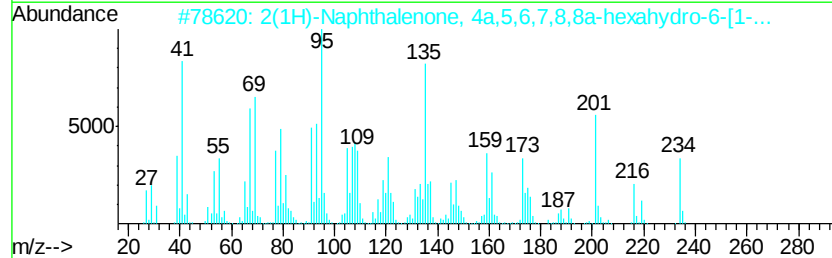
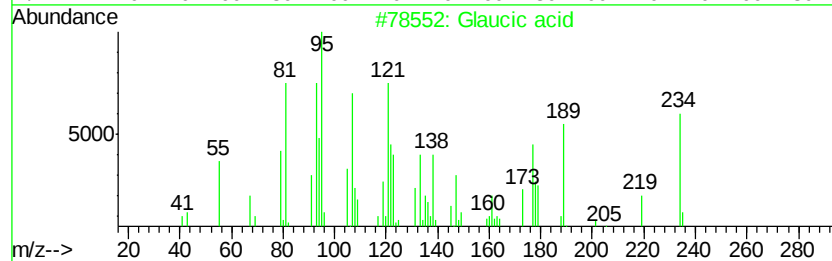
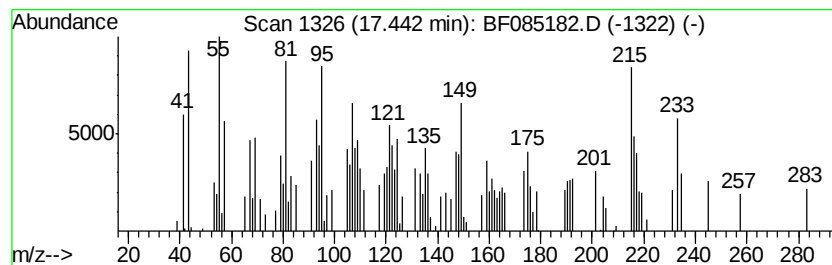
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown17.44 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.44	2.80 ng	110315	Perylene-d12	15.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Glaucic acid	234	C15H22O2	087745-30-0	44
2			2(1H)-Naphthalenone, 4a,5,6,7,8,...	234	C15H22O2	109897-57-6	25
3			Bisvertinol	498	C28H34O8	103804-06-4	22
4			Cyclopent[e]-1,3-oxazine, octahy...	216	C13H16N2O	1000138-92-2	18
5			3a,9-Dimethyldodecahydrocyclohep...	234	C16H26O	1000189-38-7	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
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ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-49(COMP)

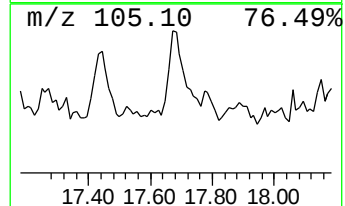
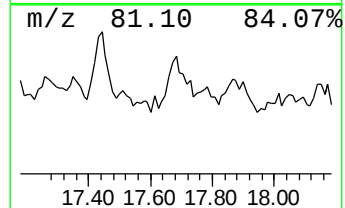
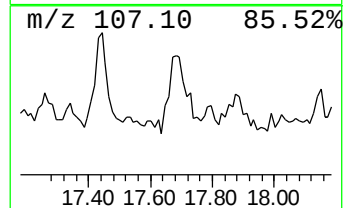
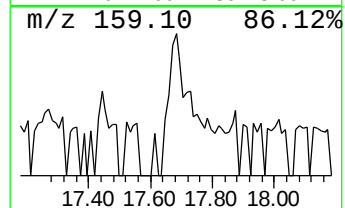
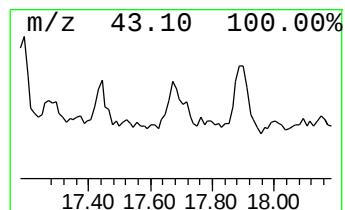
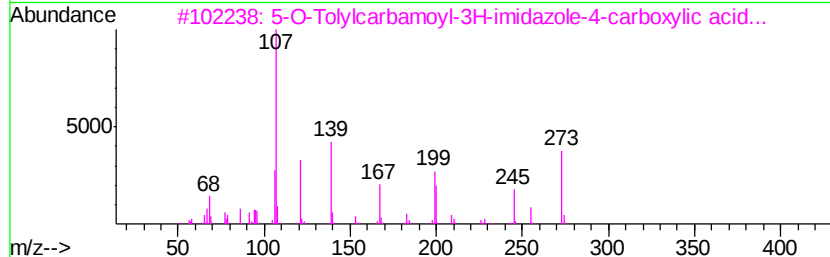
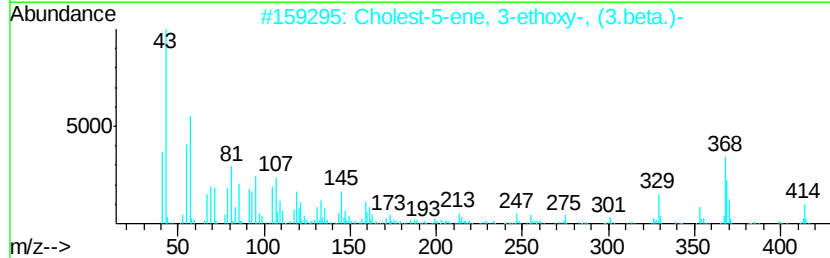
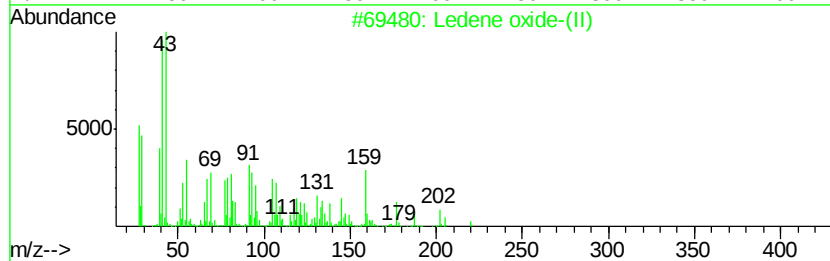
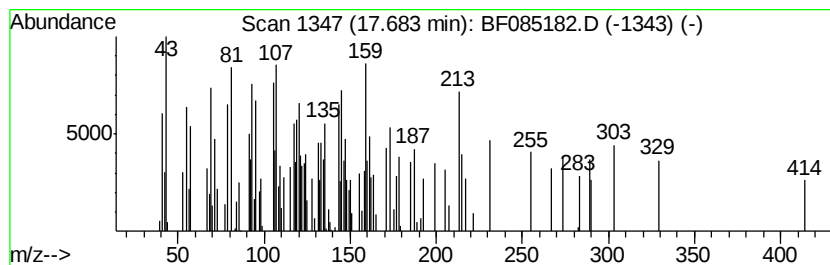
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown17.68 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.68	3.71 ng	146522	Perylene-d12	15.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ledene oxide-(II)	220	C15H24O	1000159-36-7	38
2		Cholest-5-ene, 3-ethoxy-, (3.beta...	414	C29H50O	000986-19-6	30
3		5-O-Tolylcarbamoyl-3H-imidazole-...	273	C14H15N3O3	313251-31-9	14
4		5-Benzylcarbamoyl-1H-imidazole-4...	273	C14H15N3O3	312758-83-1	14
5		5-(1-Benzoxazolyl)-2(1H)-pyrimid...	213	C11H7N3O2	349488-95-5	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF030316\
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ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-49(COMP)

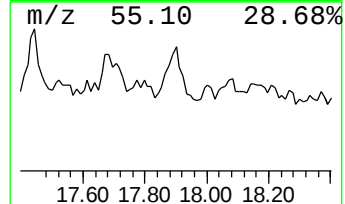
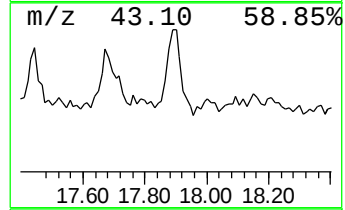
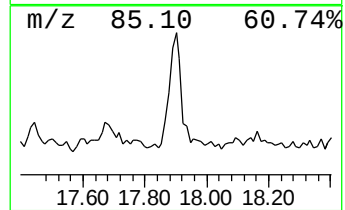
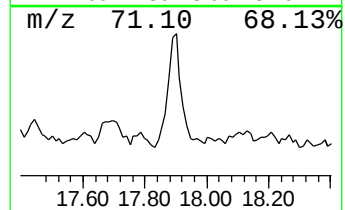
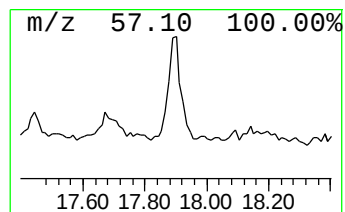
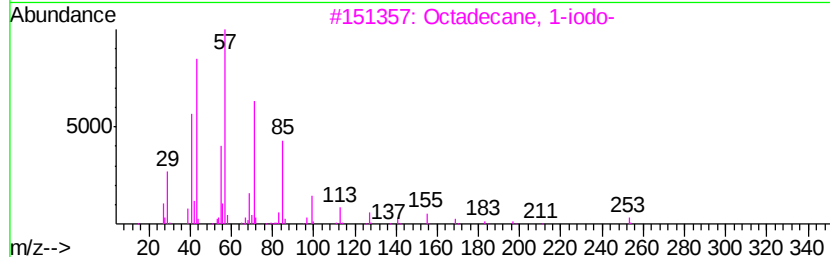
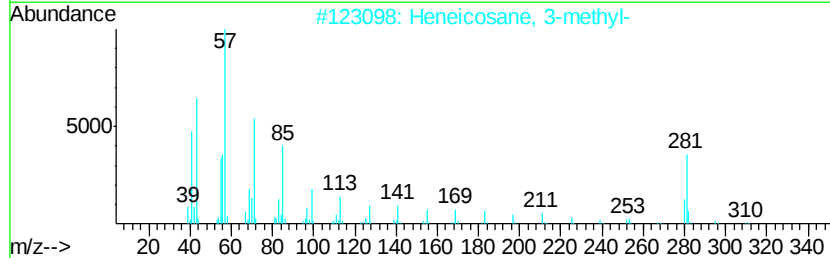
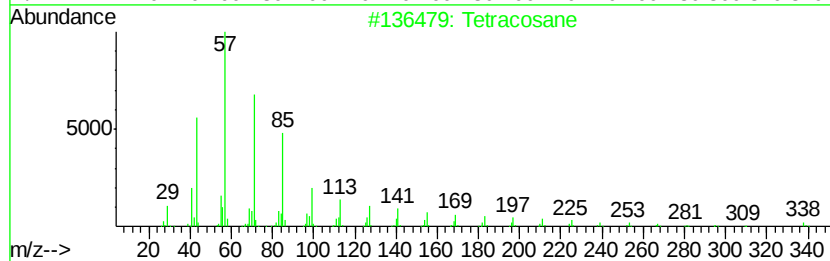
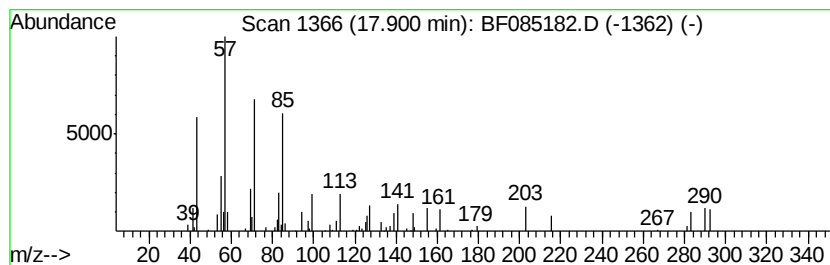
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Tetracosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.90	2.66 ng	104825	Perylene-d12	15.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	52
2			Heneicosane, 3-methyl-	310	C22H46	006418-47-9	50
3			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	43
4			Pentadecane, 2-methyl-	226	C16H34	001560-93-6	43
5			Eicosane, 10-methyl-	296	C21H44	054833-23-7	43



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 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 SB-49(COMP)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	21.0	ng	940788	1	6.97	895510	20.0
unknown6.74	6.74	85.7	ng	3836060	1	6.97	895510	20.0
Heptadecane	10.45	3.8	ng	238553	3	10.01	1270330	20.0
n-Hexadecanoic acid	12.02	6.9	ng	561519	4	11.50	1622740	20.0
2-Chloropropioni...	13.98	5.5	ng	331258	5	14.14	1197310	20.0
3-Cyclohexylpropi...	14.89	3.8	ng	147924	6	15.60	788829	20.0
Squalene	14.99	3.0	ng	116801	6	15.60	788829	20.0
Bromoacetic acid,...	16.13	2.9	ng	112257	6	15.60	788829	20.0
Eicosane	16.59	2.5	ng	96530	6	15.60	788829	20.0
unknown17.13	17.13	3.5	ng	137875	6	15.60	788829	20.0
unknown17.27	17.27	2.3	ng	92160	6	15.60	788829	20.0
unknown17.44	17.44	2.8	ng	110315	6	15.60	788829	20.0
unknown17.68	17.68	3.7	ng	146522	6	15.60	788829	20.0
Tetracosane	17.90	2.7	ng	104825	6	15.60	788829	20.0