

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020916\
 Data File : BF084967.D
 Acq On : 9 Feb 2016 22:16
 Operator : SJ/IZ
 Sample : H1329-16
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S4-B010(20-22)

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-BF020116.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	4.784	233	236	248	rVB	841069	1517049	27.57%	2.914%
2	5.230	272	275	277	rBV	4732442	5233698	95.11%	10.052%
3	6.293	365	368	370	rBV	4455474	5154384	93.67%	9.900%
4	6.407	375	378	380	rBV	5111484	5502965	100.00%	10.569%
5	6.636	396	398	399	rBV	919831	1007139	18.30%	1.934%
6	6.785	409	411	414	rVB	2941206	3562614	64.74%	6.843%
7	7.207	445	448	450	rBV	3192597	3357146	61.01%	6.448%
8	7.333	456	459	463	rVB2	92209	148273	2.69%	0.285%
9	7.676	486	489	491	rBV	64791	67416	1.23%	0.129%
10	7.916	508	510	513	rBV	1327350	1339745	24.35%	2.573%
11	9.002	602	605	607	rBV	5088554	5132994	93.28%	9.859%
12	9.402	637	640	642	rBV	1723927	1554015	28.24%	2.985%
13	9.619	652	659	661	rBV	81596	96413	1.75%	0.185%
14	9.665	661	663	665	rBV	1498263	1476763	26.84%	2.836%
15	10.111	695	702	704	rBV2	86662	132675	2.41%	0.255%
16	10.453	729	732	735	rVV	3124355	3580682	65.07%	6.877%
17	10.579	742	743	748	rVB	75433	129773	2.36%	0.249%
18	11.025	781	782	784	rBV	47122	65382	1.19%	0.126%
19	11.139	790	792	795	rBV	1165172	1477512	26.85%	2.838%
20	11.391	811	814	818	rBV	87344	118688	2.16%	0.228%
21	11.459	818	820	825	rVB2	41884	62005	1.13%	0.119%
22	11.871	854	856	860	rBV2	59204	100054	1.82%	0.192%
23	12.202	883	885	888	rBV3	40984	57902	1.05%	0.111%
24	12.351	896	898	900	rBV	154800	142565	2.59%	0.274%
25	12.568	915	917	919	rBV2	1680128	1474058	26.79%	2.831%
26	12.648	919	924	926	rVB2	180856	255705	4.65%	0.491%
27	12.728	929	931	934	rVV	2995949	4324496	78.58%	8.306%
28	12.979	950	953	954	rBV	63735	74791	1.36%	0.144%
29	13.277	976	979	981	rBV	753317	871180	15.83%	1.673%
30	13.322	981	983	984	rBV	70839	97872	1.78%	0.188%
31	13.437	990	993	994	rBV2	34831	60074	1.09%	0.115%
32	13.494	995	998	999	rBV	59373	93726	1.70%	0.180%
33	13.642	1008	1011	1017	rBV	259750	411806	7.48%	0.791%
34	13.768	1020	1022	1025	rBV	1129827	1227877	22.31%	2.358%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.962	1037	1039	1040	rBV	158268	165959	3.02%	0.319%
36	14.260	1063	1065	1072	rVB	101362	261257	4.75%	0.502%
37	14.557	1090	1091	1094	rVB	117619	104540	1.90%	0.201%
38	14.854	1114	1117	1119	rBV	95026	152619	2.77%	0.293%
39	15.128	1139	1141	1144	rBV	620304	869185	15.79%	1.669%
40	15.174	1144	1145	1154	rVB2	95949	226525	4.12%	0.435%
41	15.963	1212	1214	1216	rBV2	32759	64014	1.16%	0.123%
42	16.397	1249	1252	1256	rBV4	29910	107331	1.95%	0.206%
43	16.866	1290	1293	1299	rBV7	53897	203812	3.70%	0.391%

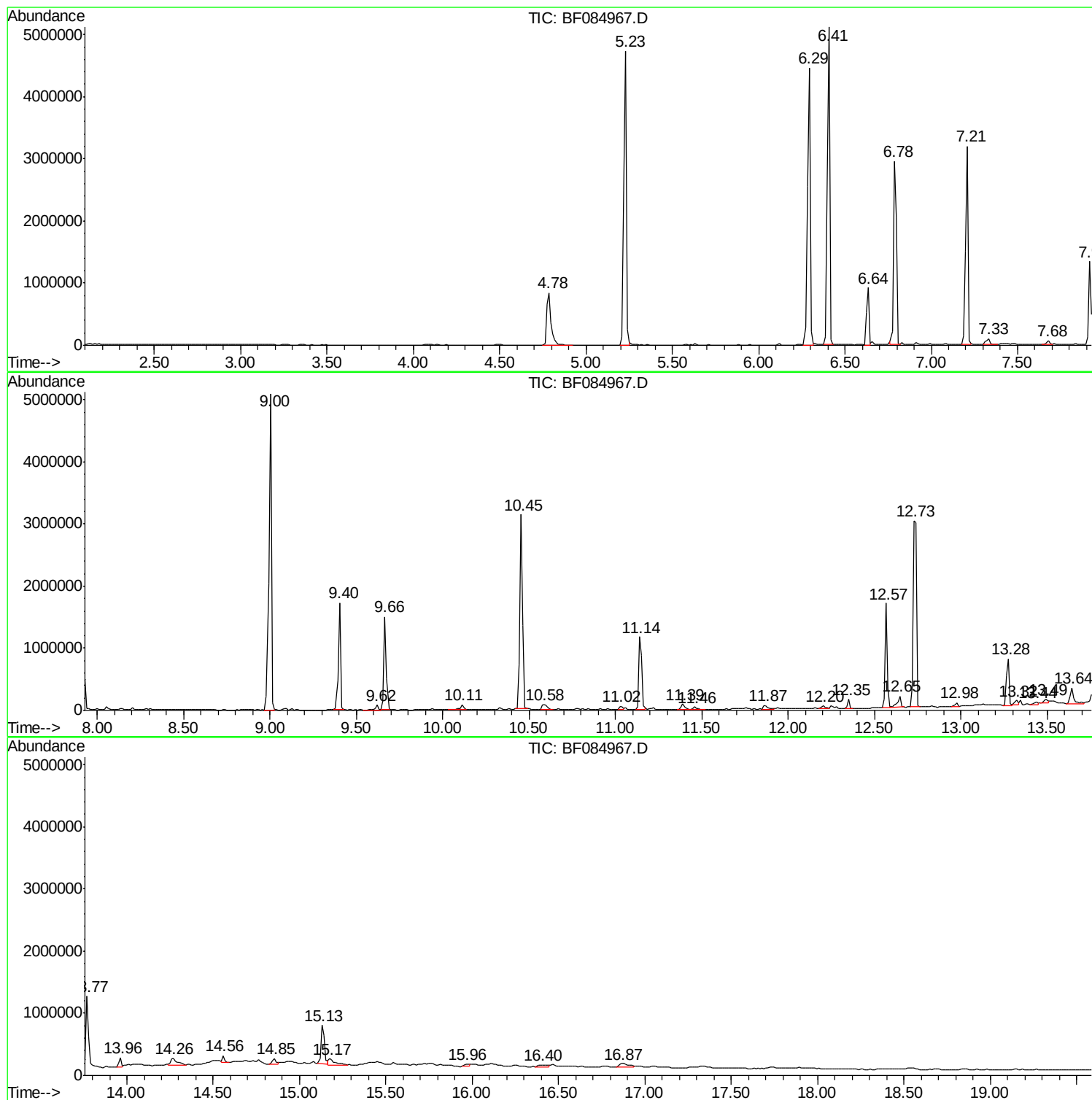
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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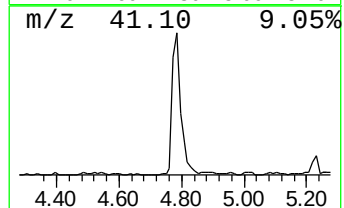
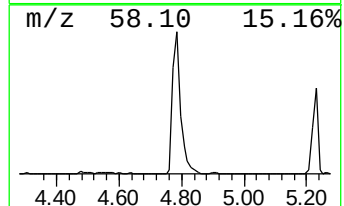
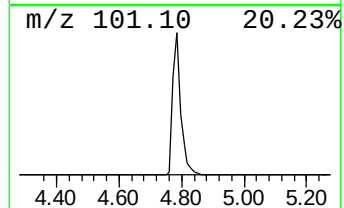
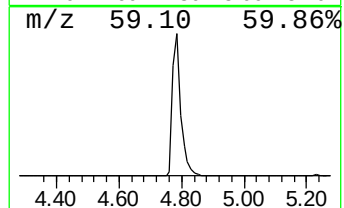
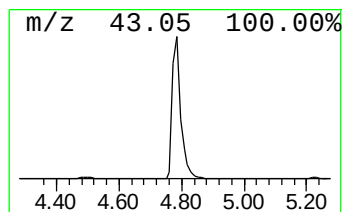
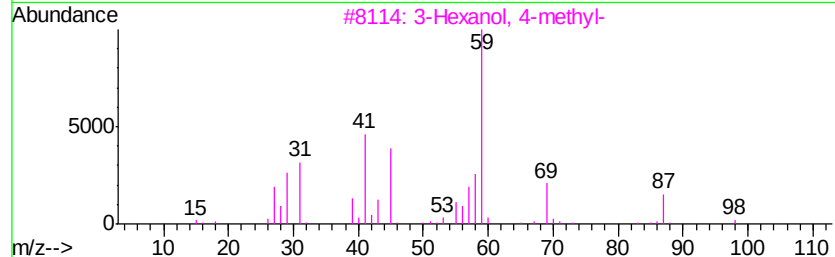
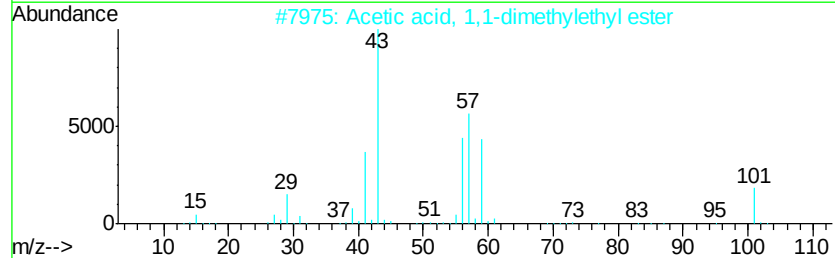
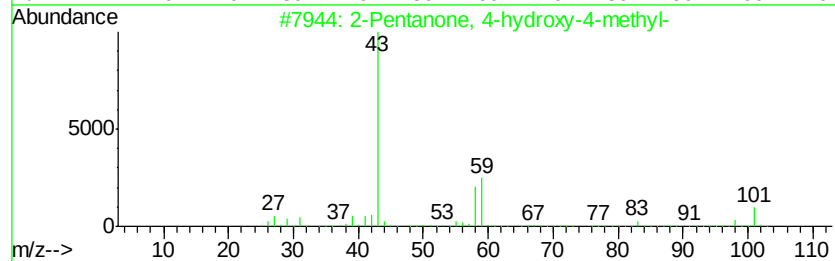
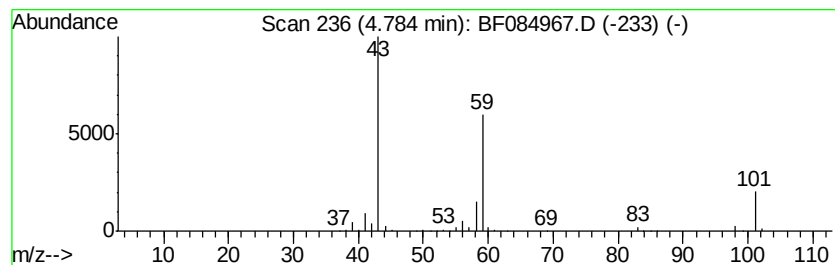
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.78	30.13 ng	1517050	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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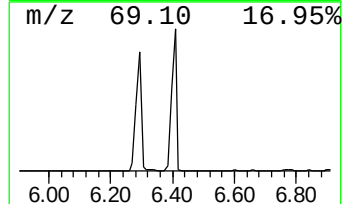
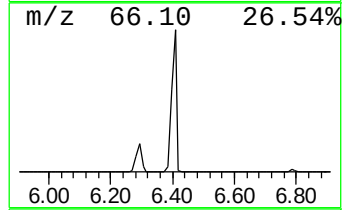
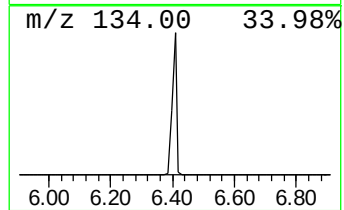
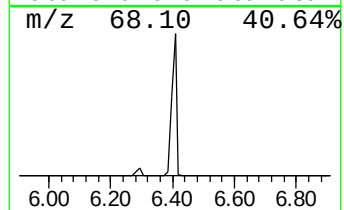
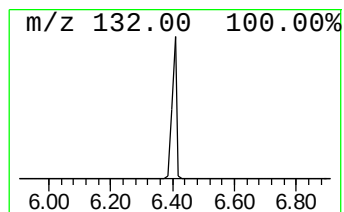
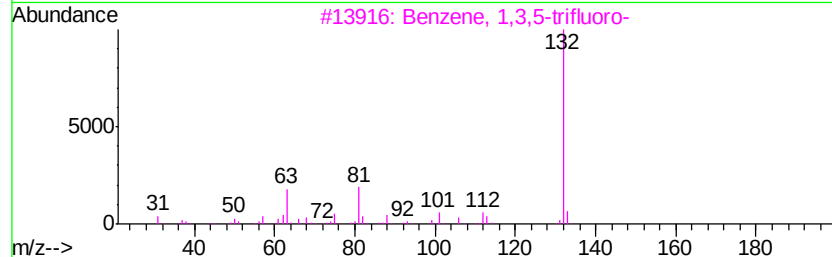
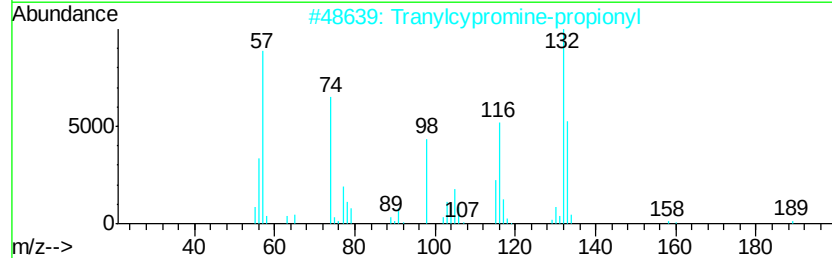
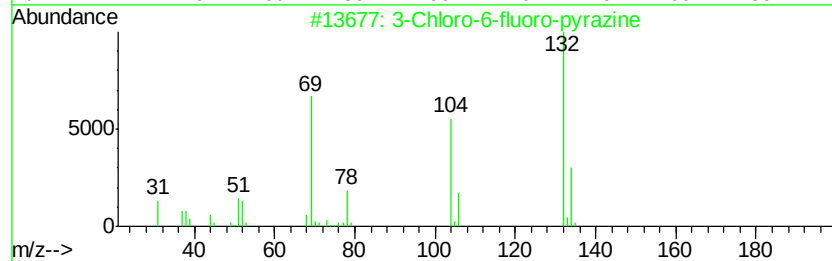
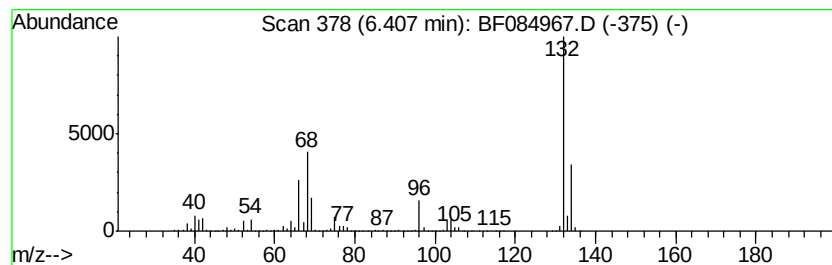
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Peak Number 2 unknown6.41 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.41	109.28 ng	5502970	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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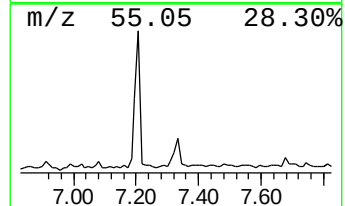
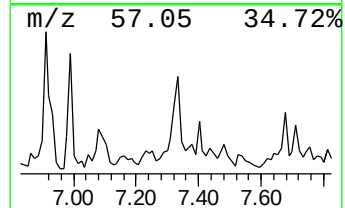
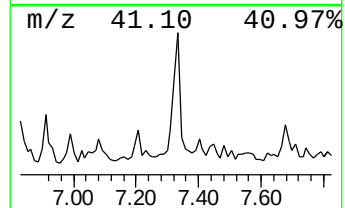
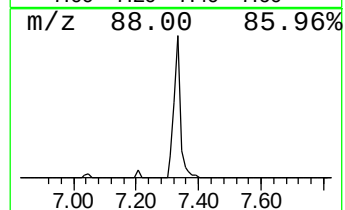
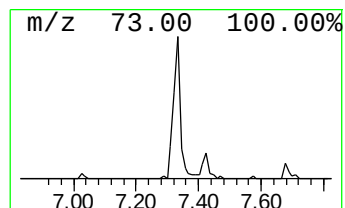
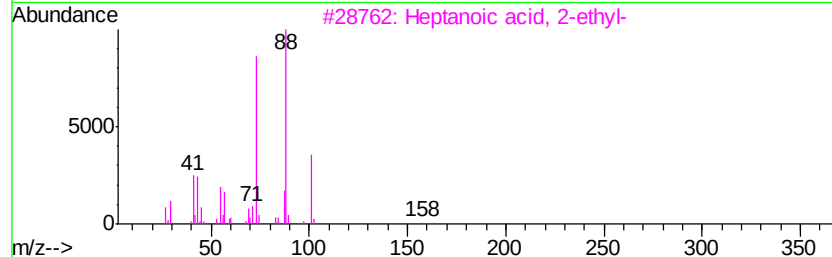
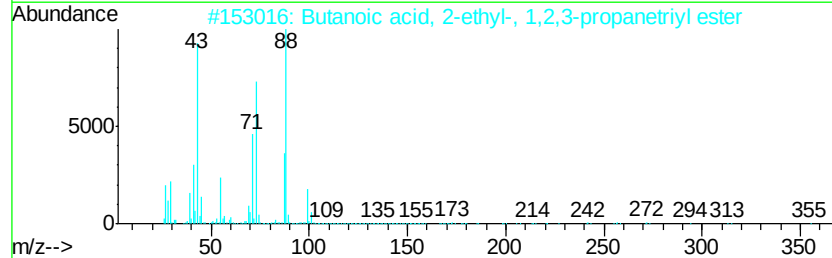
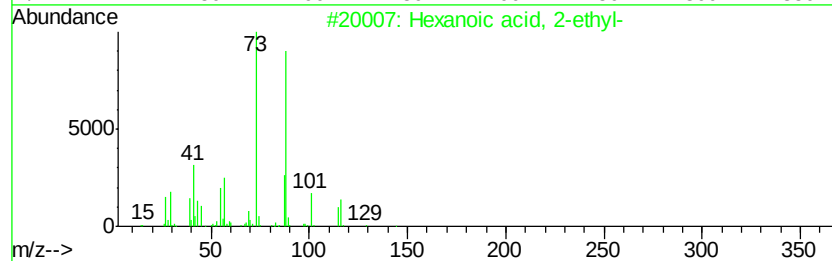
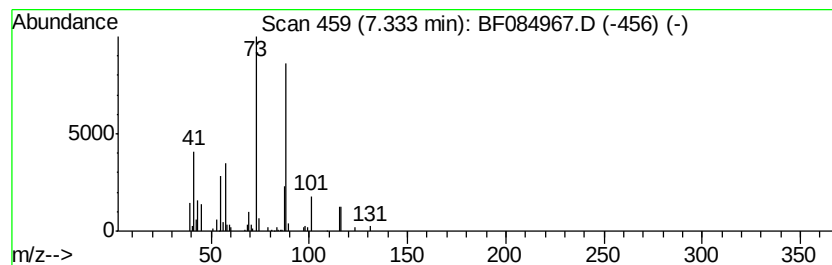
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.33	2.21 ng	148273	Naphthalene-d8	7.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2		Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	50
3		Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	42
4		1,4-Dioxane	88	C4H8O2	000123-91-1	38
5		Acetic acid, diethyl-	116	C6H12O2	000088-09-5	36



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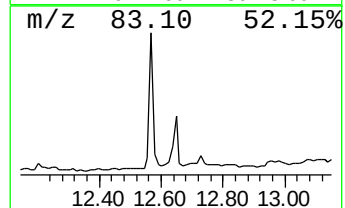
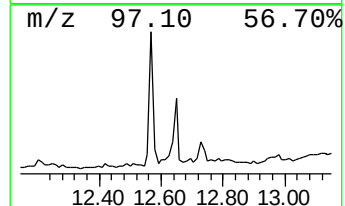
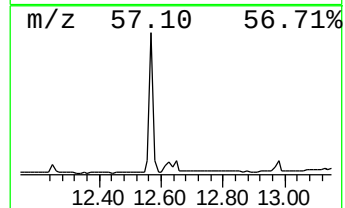
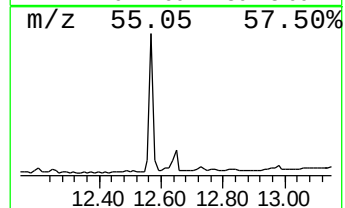
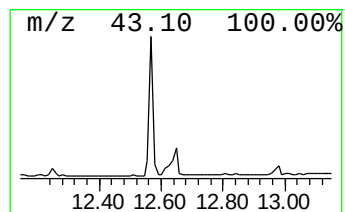
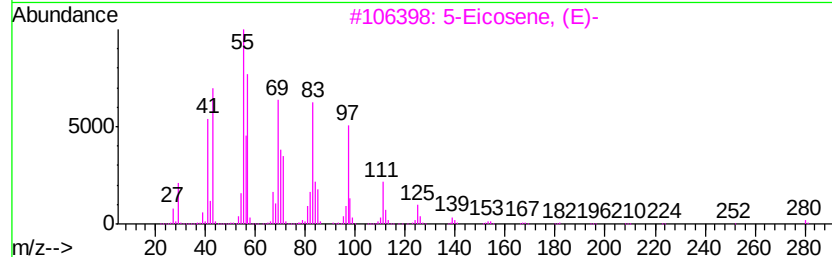
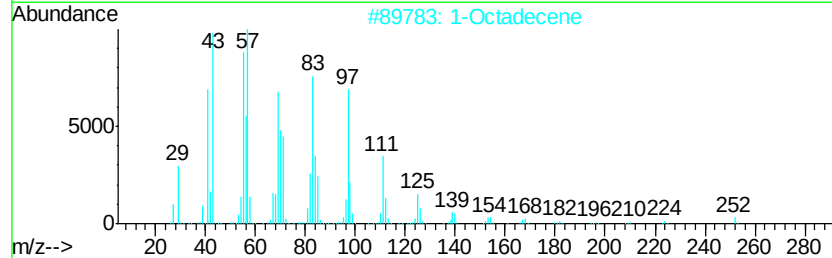
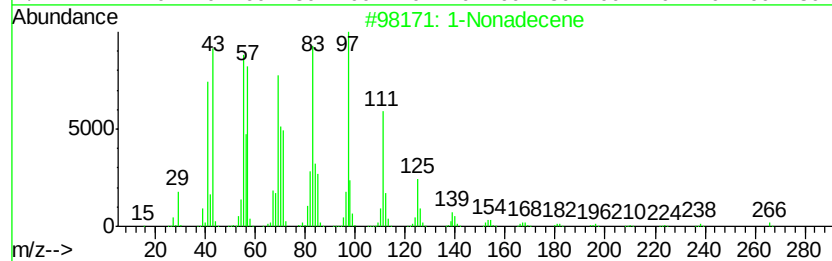
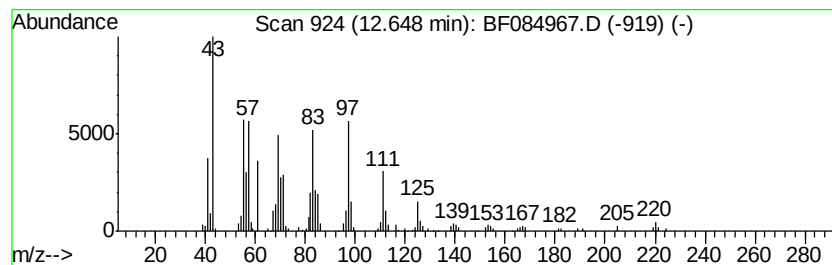
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 1-Nonadecene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	4.16 ng	255705	Chrysene-d12	13.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Nonadecene		266	C19H38	018435-45-5	93
2	1-Octadecene		252	C18H36	000112-88-9	93
3	5-Eicosene, (E)-		280	C20H40	074685-30-6	83
4	Dichloroacetic acid, tetradecyl ...		324	C16H30Cl2O2	083005-02-1	68
5	Trichloroacetic acid, tetradecyl...		358	C16H29Cl3O2	074339-52-9	68



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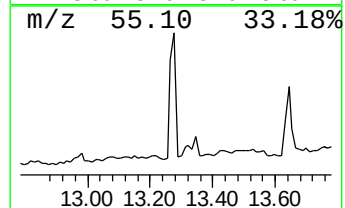
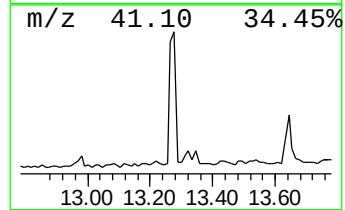
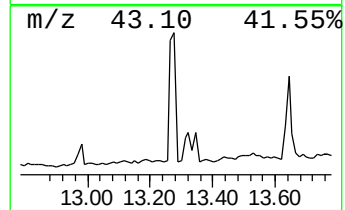
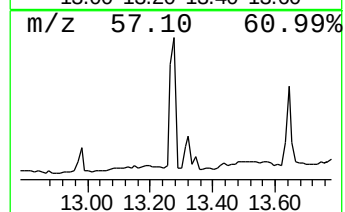
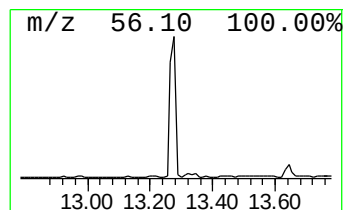
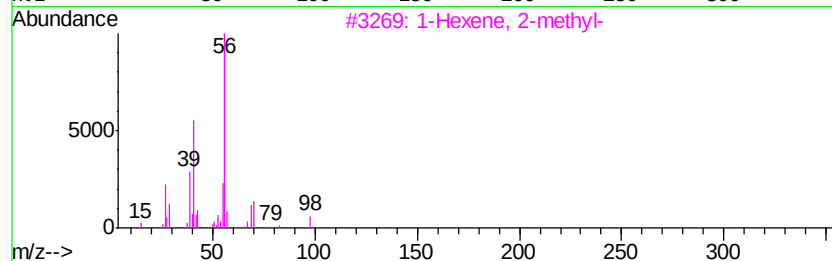
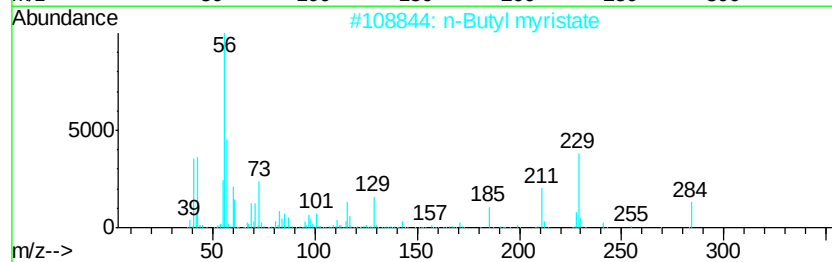
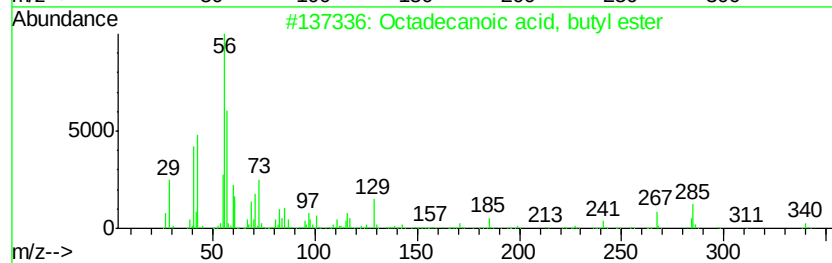
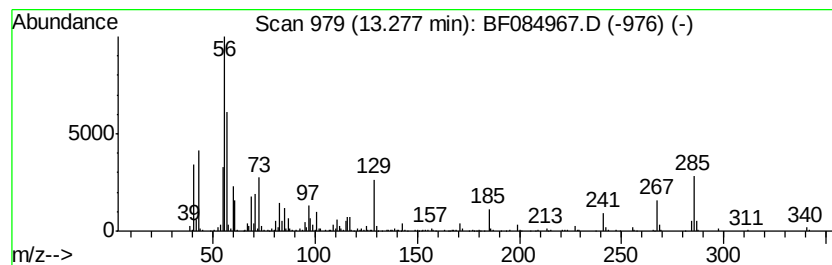
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Octadecanoic acid, butyl ester Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.28	14.19 ng	871180	Chrysene-d12	13.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	94
2		n-Butyl myristate	284	C18H36O2	000110-36-1	53
3		1-Hexene, 2-methyl-	98	C7H14	006094-02-6	27
4		1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	27
5		Cyclohexane, (1,1-dimethylethyl)-	140	C10H20	003178-22-1	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020916\
Data File : BF084967.D
Acq On : 9 Feb 2016 22:16
Operator : SJ/IZ
Sample : H1329-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S4-B010(20-22)

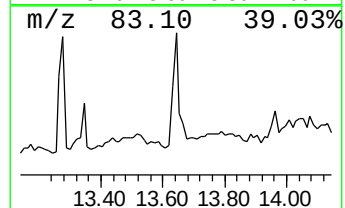
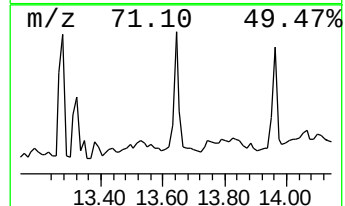
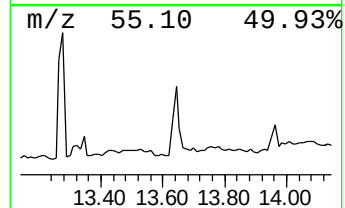
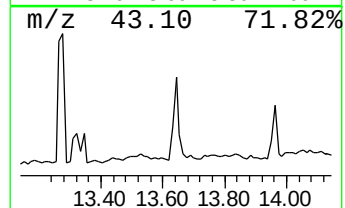
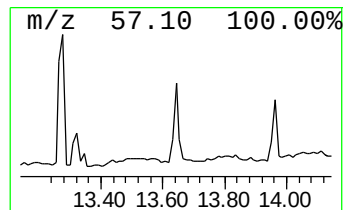
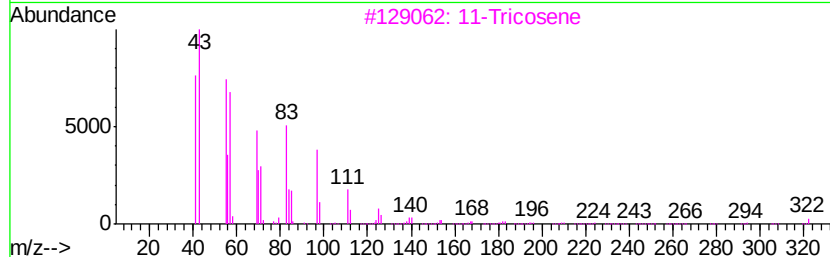
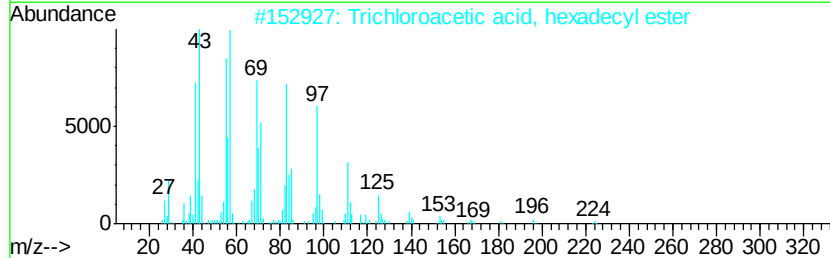
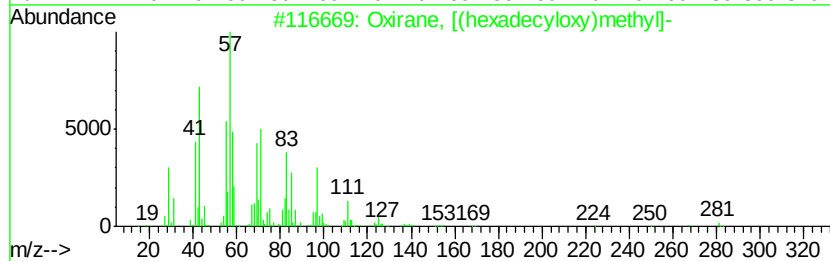
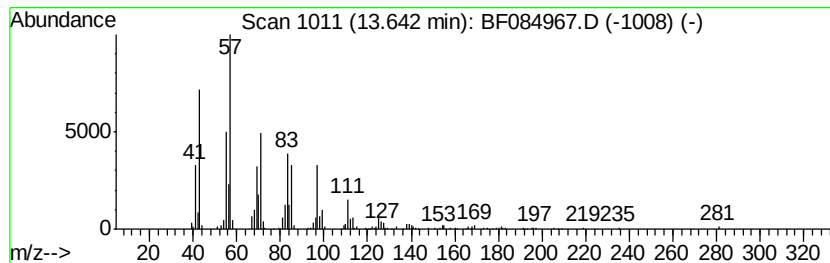
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Oxirane, [(hexadecyloxy)met... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	6.71 ng	411806	Chrysene-d12	13.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, [(hexadecyloxy)methyl]-	298	C19H38O2	015965-99-8	80
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	62
3		11-Tricosene	322	C23H46	052078-56-5	62
4		Decane	142	C10H22	000124-18-5	60
5		Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020916\
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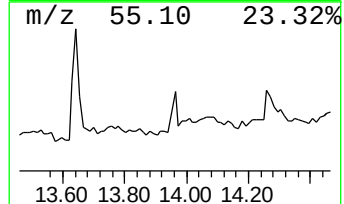
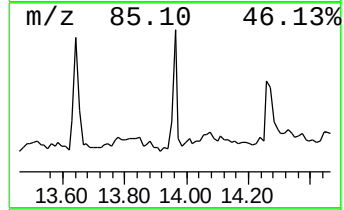
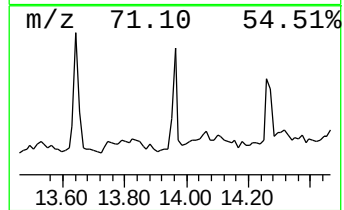
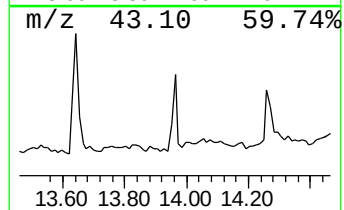
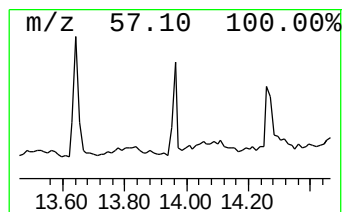
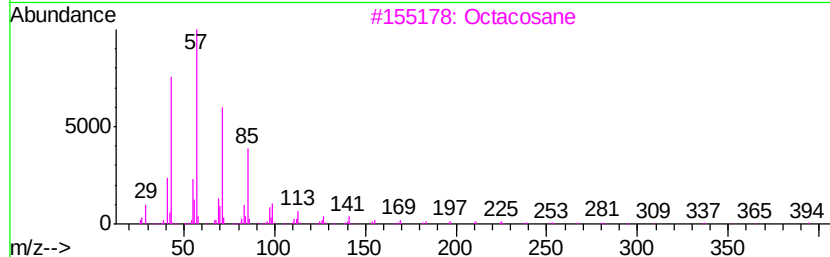
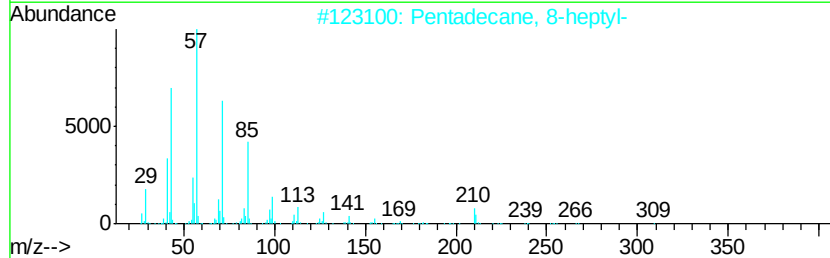
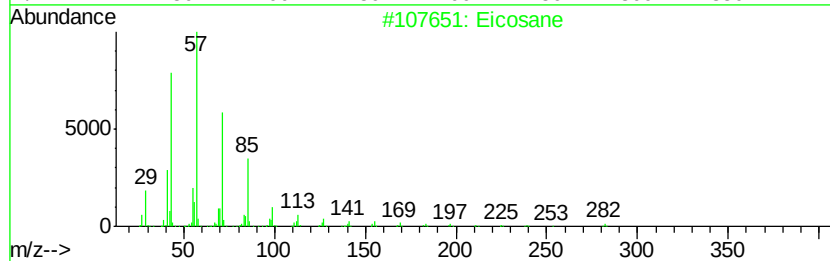
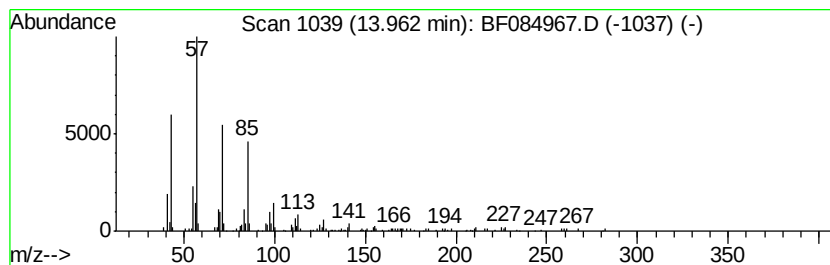
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	2.70 ng	165959	Chrysene-d12	13.77
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane		282 C20H42	000112-95-8 96
2	Pentadecane, 8-heptyl-		310 C22H46	071005-15-7 87
3	Octacosane		394 C28H58	000630-02-4 76
4	Heneicosane		296 C21H44	000629-94-7 74
5	Hexadecane		226 C16H34	000544-76-3 72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020916\
Data File : BF084967.D
Acq On : 9 Feb 2016 22:16
Operator : SJ/IZ
Sample : H1329-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

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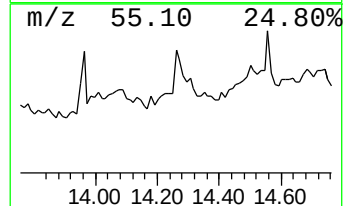
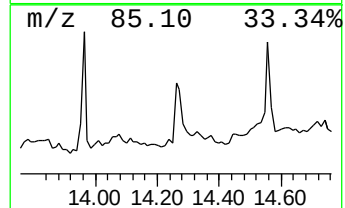
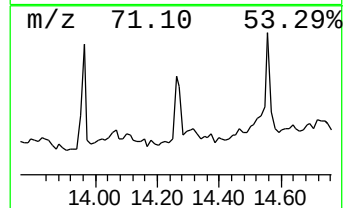
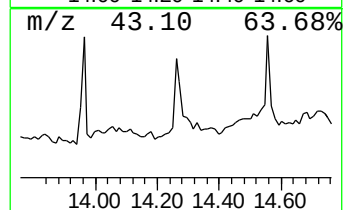
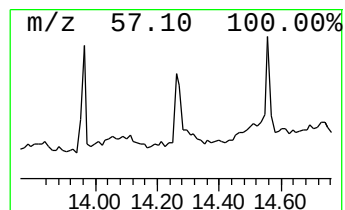
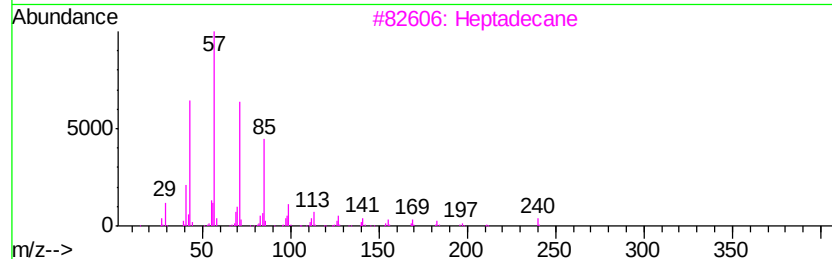
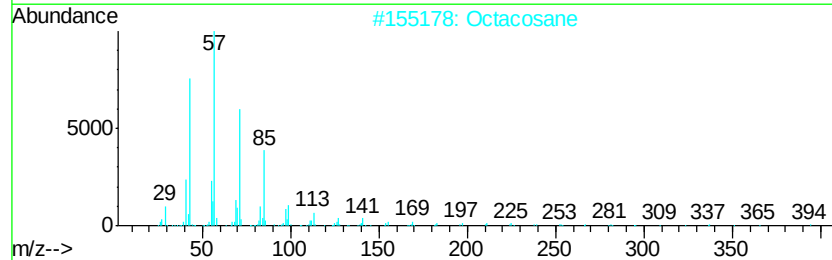
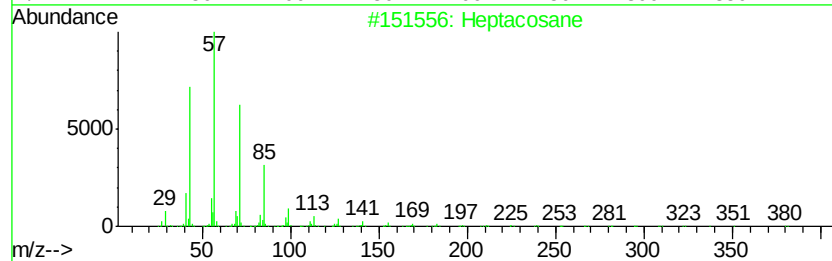
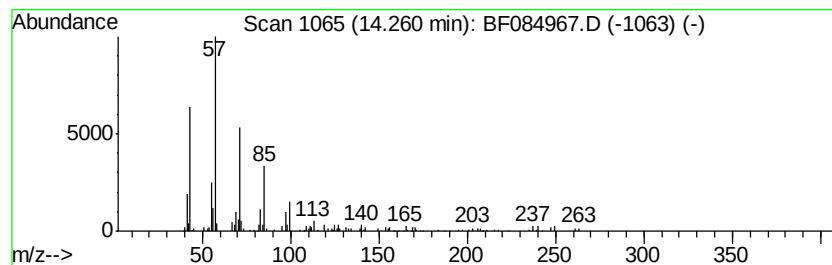
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	4.26 ng	261257	Chrysene-d12	13.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane		380	C27H56	000593-49-7	72
2	Octacosane		394	C28H58	000630-02-4	68
3	Heptadecane		240	C17H36	000629-78-7	62
4	Eicosane, 10-methyl-		296	C21H44	054833-23-7	59
5	Tetracosane		338	C24H50	000646-31-1	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020916\
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Acq On : 9 Feb 2016 22:16
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Sample : H1329-16
Misc :
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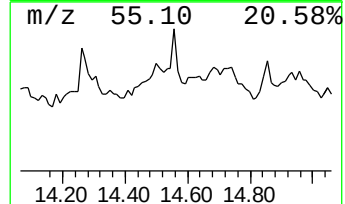
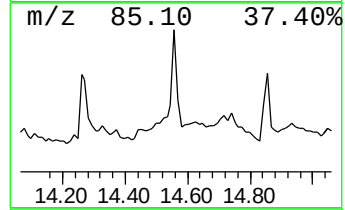
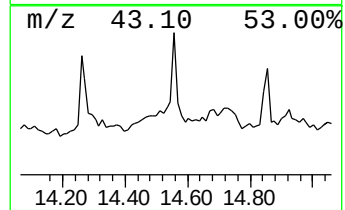
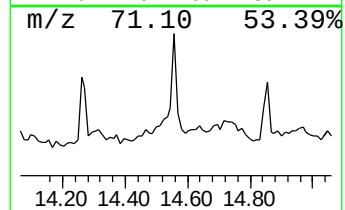
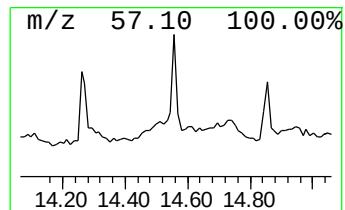
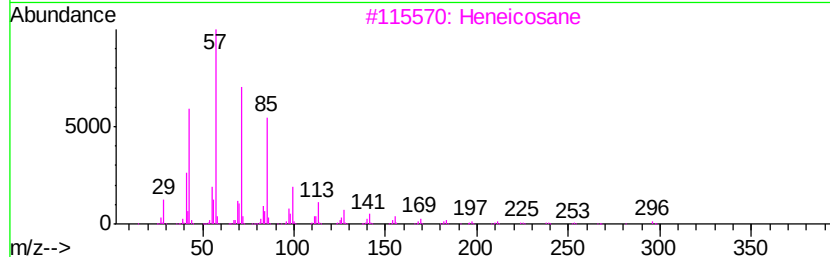
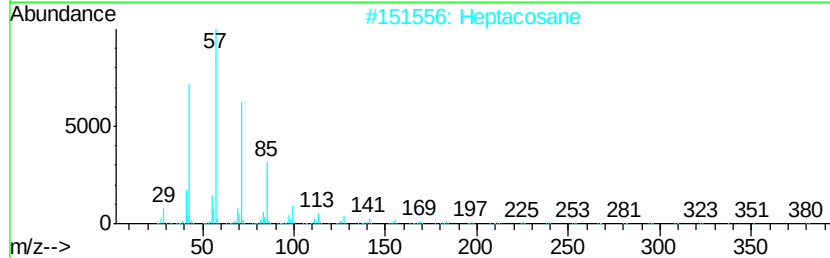
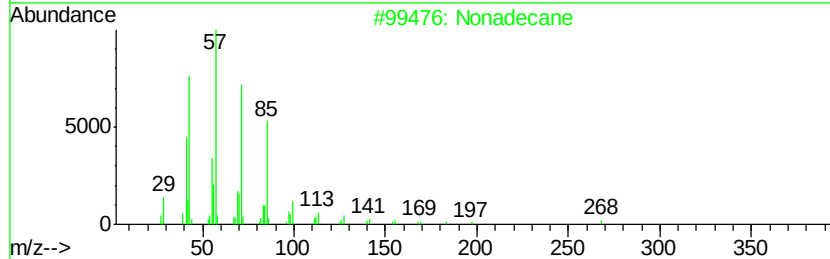
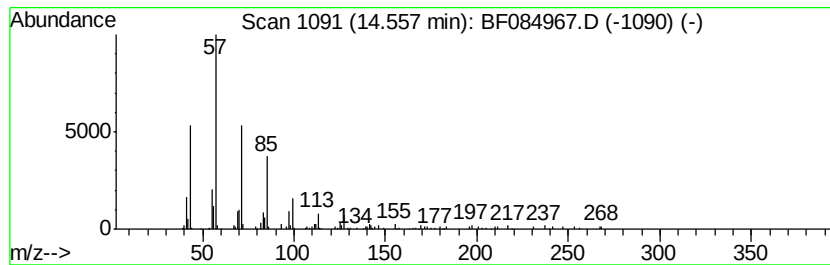
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Nonadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.56	2.41 ng	104540	Perylene-d12	15.13	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	81
2	Heptacosane	380	C27H56	000593-49-7	80
3	Heneicosane	296	C21H44	000629-94-7	80
4	Eicosane	282	C20H42	000112-95-8	80
5	Heptadecane	240	C17H36	000629-78-7	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020916\
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ClientSampleId :
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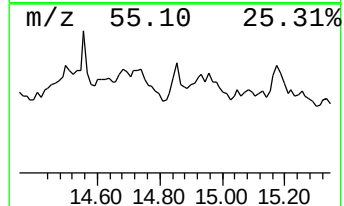
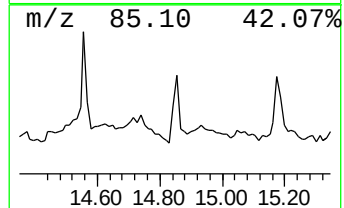
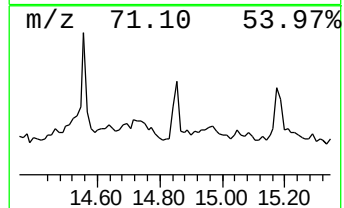
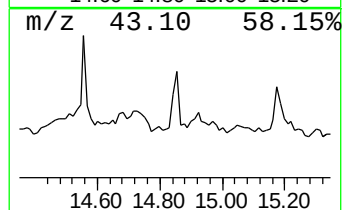
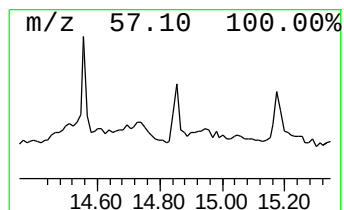
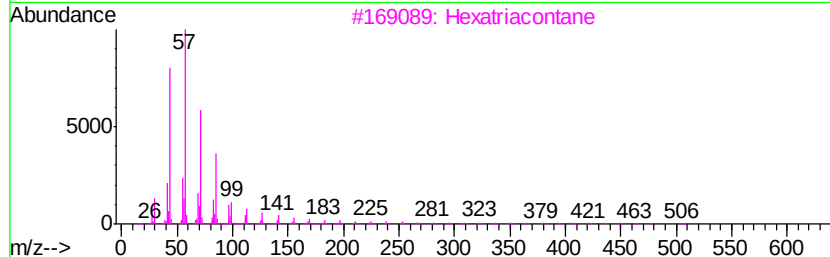
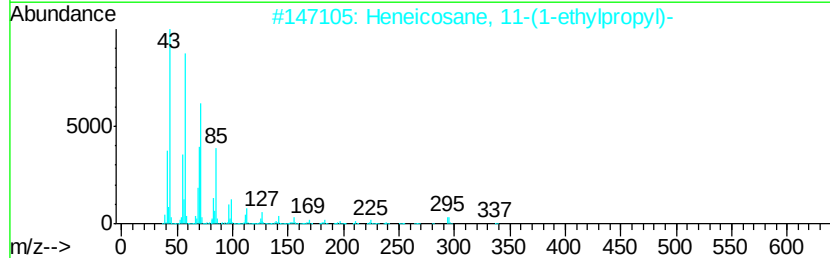
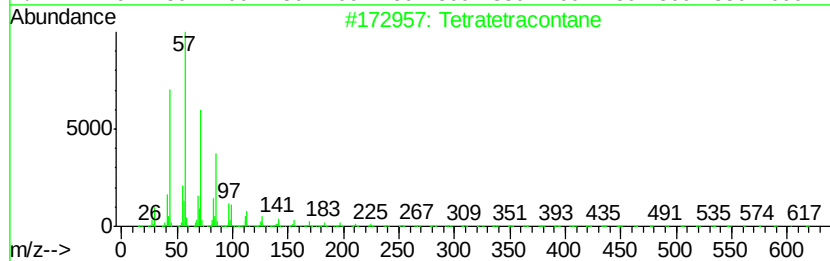
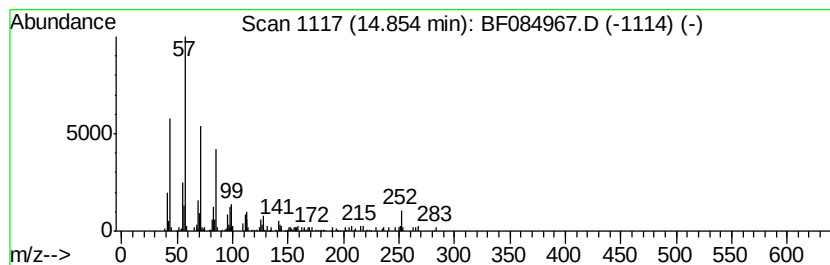
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Tetratetracontane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	3.51 ng	152619	Perylene-d12	15.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	87
2			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	83
3			Hexatriacontane	507	C36H74	000630-06-8	83
4			Hentriacontane	437	C31H64	000630-04-6	80
5			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020916\
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Acq On : 9 Feb 2016 22:16
Operator : SJ/IZ
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Misc :
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ClientSampleId :
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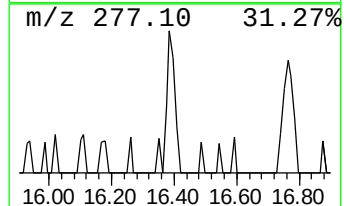
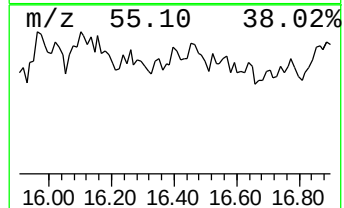
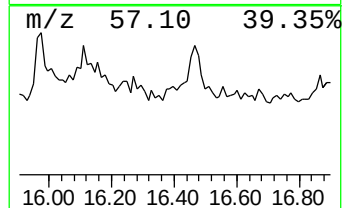
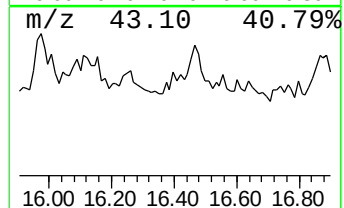
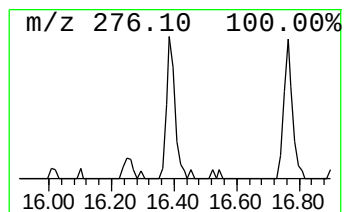
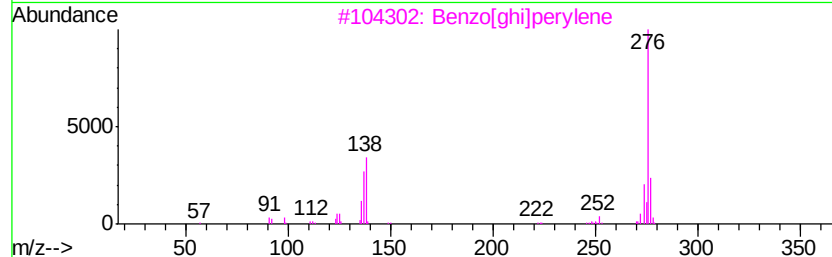
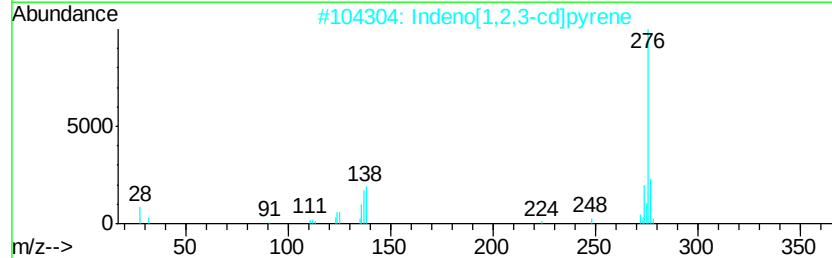
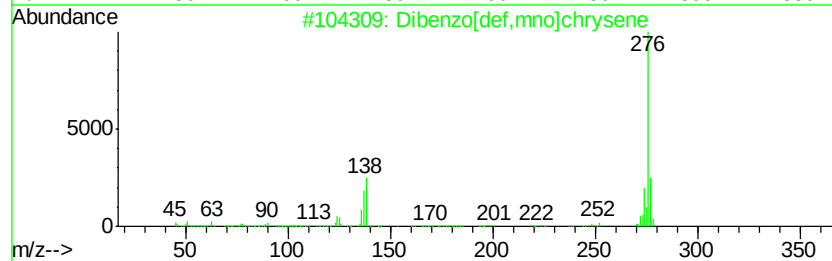
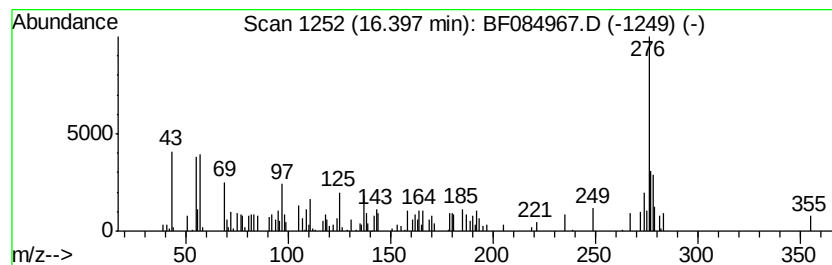
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown16.40 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	2.47 ng	107331	Perylene-d12	15.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	49
2		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	49
3		Benzo[ghi]perylene	276	C22H12	000191-24-2	43
4		9,10-Anthracenedione, 1,8-dichloro-	276	C14H6Cl2O2	000082-43-9	43
5		Acethydrazide, N2-(4-iodophenyl)-	276	C8H9IN2O	014579-96-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020916\
Data File : BF084967.D
Acq On : 9 Feb 2016 22:16
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Misc :
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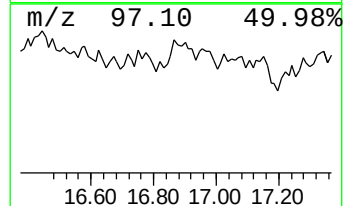
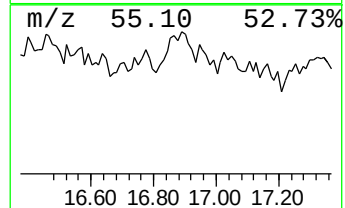
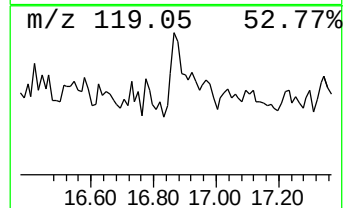
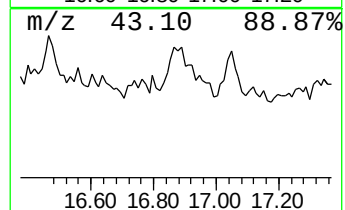
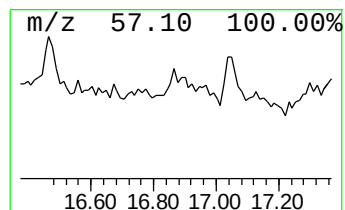
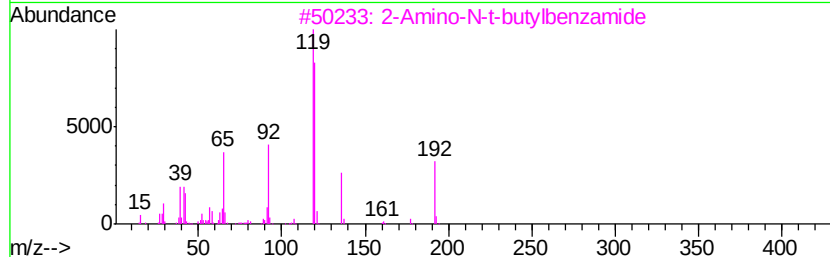
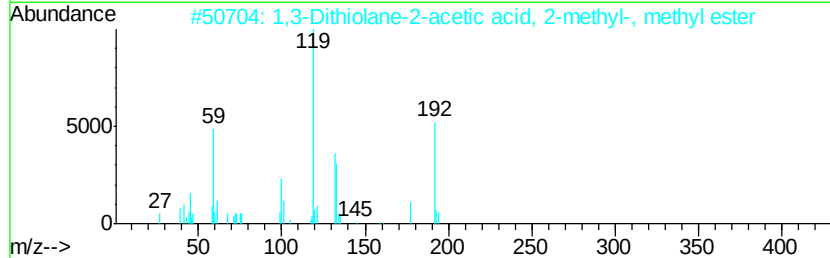
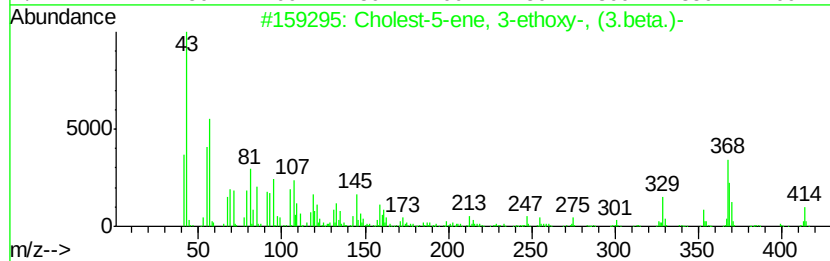
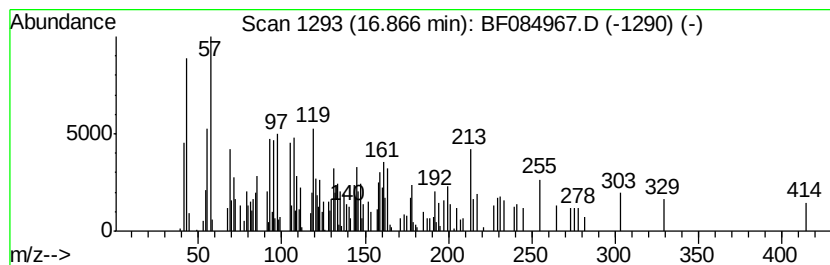
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown16.87 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.87	4.69 ng	203812	Perylene-d12	15.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cholest-5-ene, 3-ethoxy-, (3.beta...	414	C29H50O	000986-19-6	30
2		1,3-Dithiolane-2-acetic acid, 2-...	192	C7H12O2S2	065726-45-6	10
3		2-Amino-N-t-butylbenzamide	192	C11H16N2O	001203-89-0	9
4		2,3-Dihydrobenzofuran-3-carboxyl...	192	C11H12O3	1000194-68-1	9
5		5-Isopropyl-2,2,7a-trimethylhexa...	228	C13H24O3	075316-52-8	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020916\
Data File : BF084967.D
Acq On : 9 Feb 2016 22:16
Operator : SJ/IJ
Sample : H1329-16
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S4-B010(20-22)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.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...	4.78	30.1	ng	1517050	1	6.64	1007140	20.0
unknown6.41	6.41	109.3	ng	5502970	1	6.64	1007140	20.0
Hexanoic acid, 2-...	7.33	2.2	ng	148273	2	7.92	1339750	20.0
1-Nonadecene	12.65	4.2	ng	255705	5	13.77	1227880	20.0
Octadecanoic acid...	13.28	14.2	ng	871180	5	13.77	1227880	20.0
Oxirane, [(hexade...	13.64	6.7	ng	411806	5	13.77	1227880	20.0
Eicosane	13.96	2.7	ng	165959	5	13.77	1227880	20.0
Heptacosane	14.26	4.3	ng	261257	5	13.77	1227880	20.0
Nonadecane	14.56	2.4	ng	104540	6	15.13	869185	20.0
Tetratetracontane	14.85	3.5	ng	152619	6	15.13	869185	20.0
unknown16.40	16.40	2.5	ng	107331	6	15.13	869185	20.0
unknown16.87	16.87	4.7	ng	203812	6	15.13	869185	20.0