

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011317\
 Data File : BF092264.D
 Acq On : 14 Jan 2017 1:42
 Operator : UM/SJ
 Sample : I1031-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-0027

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-BF122116.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.312	43	46	52	rVB	60596	103527	1.83%	0.201%
2	4.701	252	255	265	rBV	1310482	2334251	41.37%	4.541%
3	5.170	294	296	299	rBV	4186024	4707371	83.44%	9.157%
4	6.244	387	390	392	rBV	4988387	5641951	100.00%	10.975%
5	6.347	397	399	401	rBV	5316515	4962215	87.95%	9.652%
6	6.576	417	419	421	rBV	1016427	1176968	20.86%	2.289%
7	6.724	430	432	435	rBV	3522698	3991886	70.75%	7.765%
8	7.147	466	469	471	rBV	2978175	3126859	55.42%	6.082%
9	7.296	479	482	486	rBV	61168	95809	1.70%	0.186%
10	7.856	529	531	534	rBV	1462900	1636703	29.01%	3.184%
11	8.942	623	626	628	rBV	4489740	4623179	81.94%	8.993%
12	9.342	658	661	664	rBV	1377097	1230306	21.81%	2.393%
13	9.605	682	684	687	rVB	1439588	1528678	27.09%	2.974%
14	10.062	719	724	725	rBV	77078	102549	1.82%	0.199%
15	10.393	751	753	756	rBV2	2098365	2936779	52.05%	5.713%
16	10.530	763	765	769	rVB	121553	145848	2.59%	0.284%
17	10.976	802	804	805	rBV2	73528	71634	1.27%	0.139%
18	11.079	811	813	816	rBV	1197465	1549907	27.47%	3.015%
19	11.331	833	835	839	rVB	91492	115622	2.05%	0.225%
20	11.639	860	862	866	rBV2	321112	314355	5.57%	0.611%
21	12.291	917	919	921	rBV	81080	69780	1.24%	0.136%
22	12.519	936	939	941	rBV	66210	86057	1.53%	0.167%
23	12.565	941	943	949	rVB	48350	59779	1.06%	0.116%
24	12.668	949	952	955	rBV	3239205	3564073	63.17%	6.933%
25	12.919	970	974	975	rBV2	68100	95500	1.69%	0.186%
26	13.262	1001	1004	1007	rVB	35502	64951	1.15%	0.126%
27	13.582	1029	1032	1041	rBV	300438	473700	8.40%	0.921%
28	13.708	1041	1043	1046	rBV	1297713	1286858	22.81%	2.503%
29	14.211	1084	1087	1091	rBV	99361	159520	2.83%	0.310%
30	14.554	1115	1117	1121	rVB	115163	140617	2.49%	0.274%
31	14.622	1121	1123	1126	rBV	622717	588615	10.43%	1.145%
32	14.805	1135	1139	1148	rBV	602007	1046598	18.55%	2.036%
33	15.068	1159	1162	1164	rBV	698158	910932	16.15%	1.772%
34	15.274	1177	1180	1183	rVB	62054	81801	1.45%	0.159%

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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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.468	1194	1197	1199	rBV	93163	144547	2.56%	0.281%
36	15.514	1199	1201	1205	rVV4	47232	99293	1.76%	0.193%
37	15.880	1230	1233	1239	rVB2	28296	59743	1.06%	0.116%
38	16.120	1250	1254	1257	rBV3	35140	69089	1.22%	0.134%
39	16.291	1266	1269	1272	rBV3	44574	103269	1.83%	0.201%
40	16.405	1272	1279	1284	rVV2	505009	1179113	20.90%	2.294%
41	16.497	1284	1287	1288	rVV	59635	110376	1.96%	0.215%
42	16.542	1288	1291	1295	rVB2	80787	185041	3.28%	0.360%
43	16.760	1306	1310	1316	rBV3	68700	183685	3.26%	0.357%
44	17.023	1329	1333	1335	rBV4	23612	62070	1.10%	0.121%
45	17.583	1378	1382	1388	rVB4	38249	95817	1.70%	0.186%
46	18.051	1419	1423	1428	rVB6	31450	92182	1.63%	0.179%

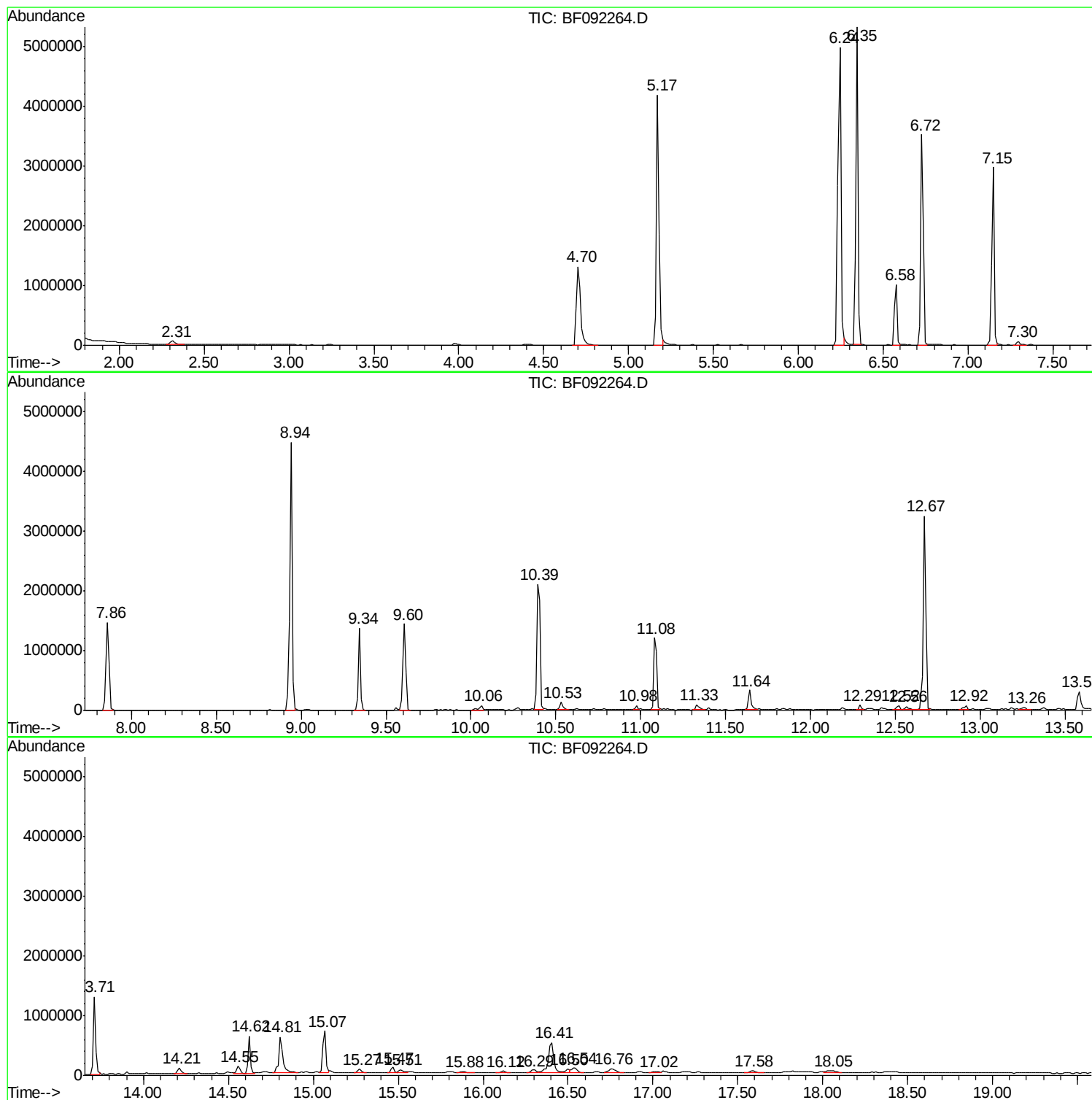
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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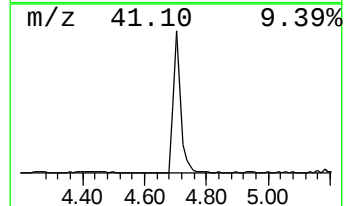
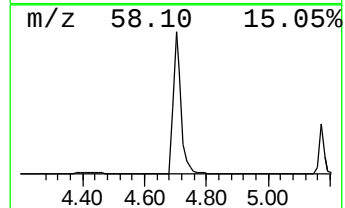
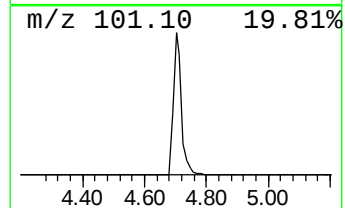
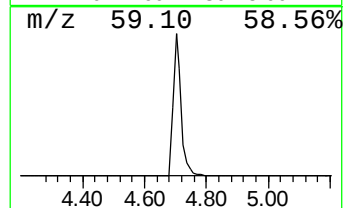
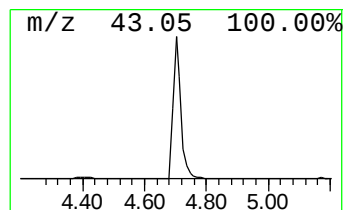
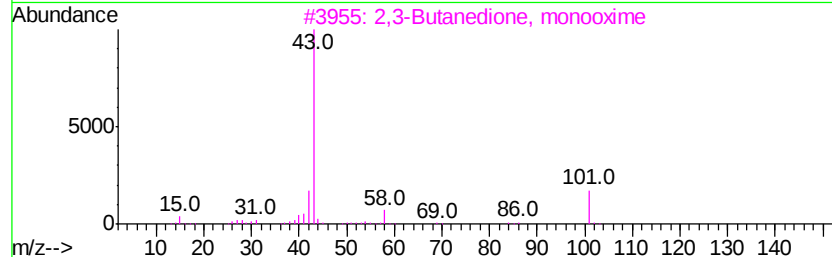
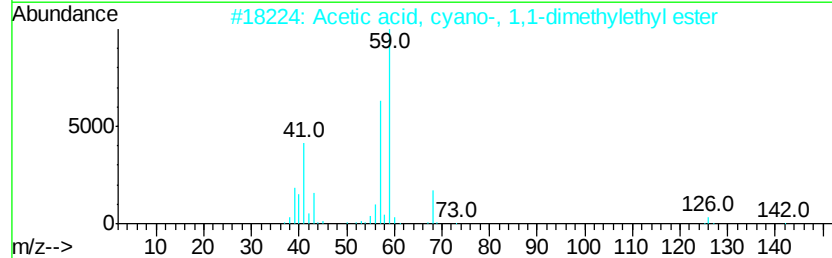
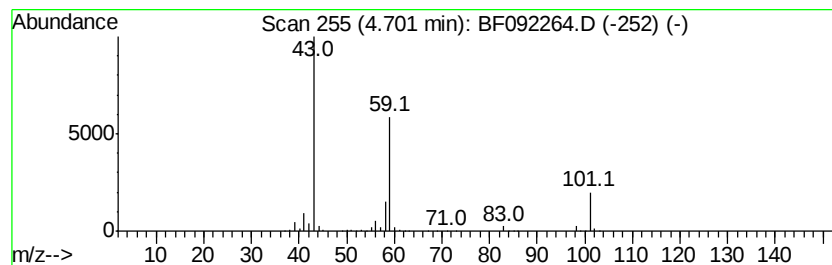
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.70	39.67 ng	2334250	1,4-Dichlorobenzene-d4	6.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	9



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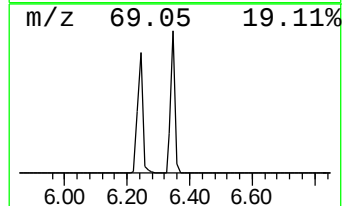
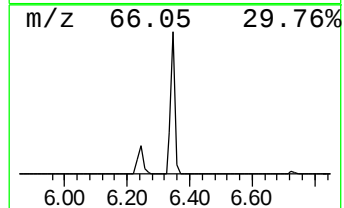
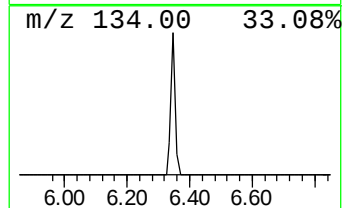
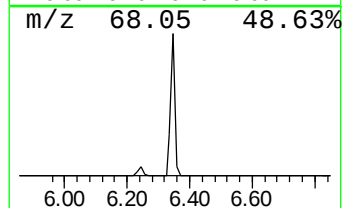
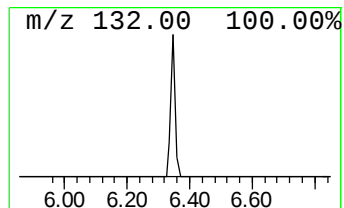
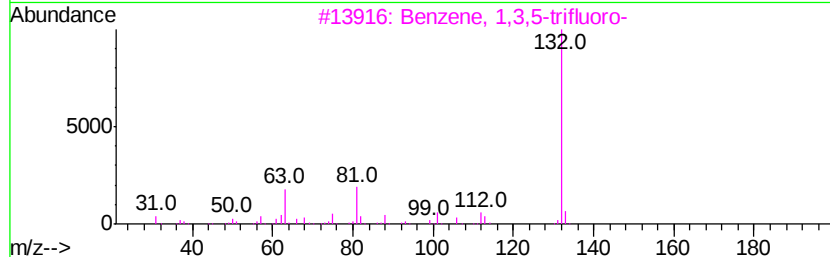
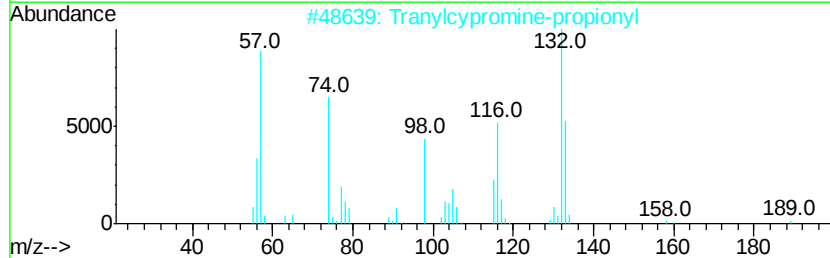
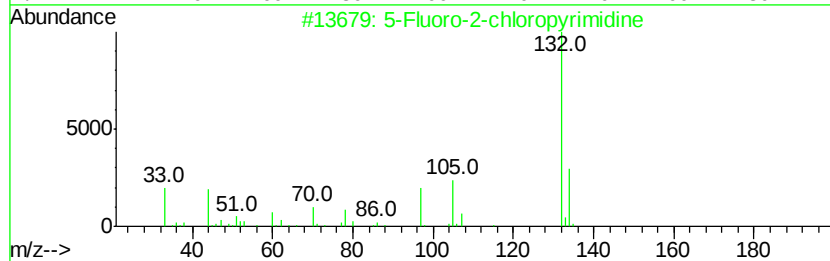
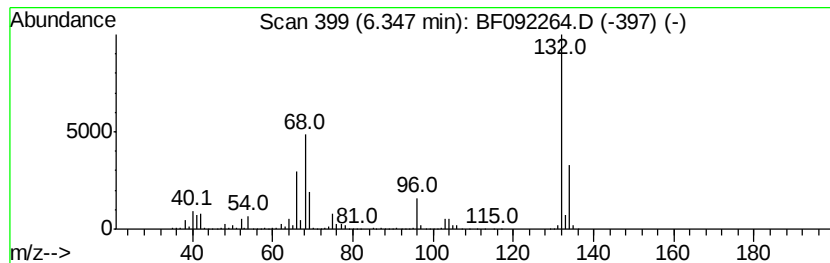
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.35 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	84.32 ng	4962220	1,4-Dichlorobenzene-d4	6.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
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		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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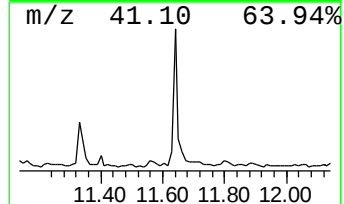
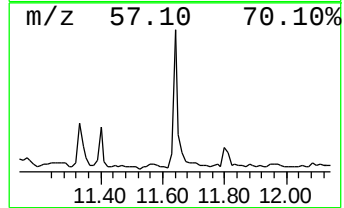
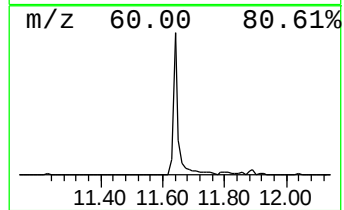
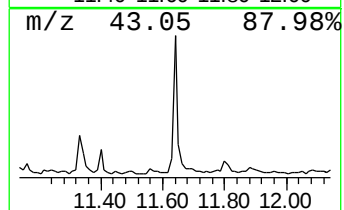
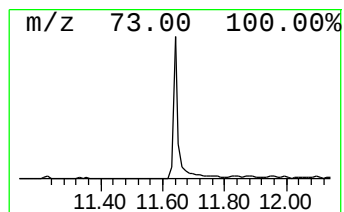
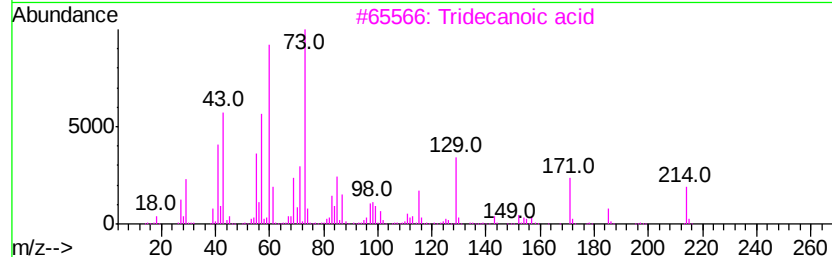
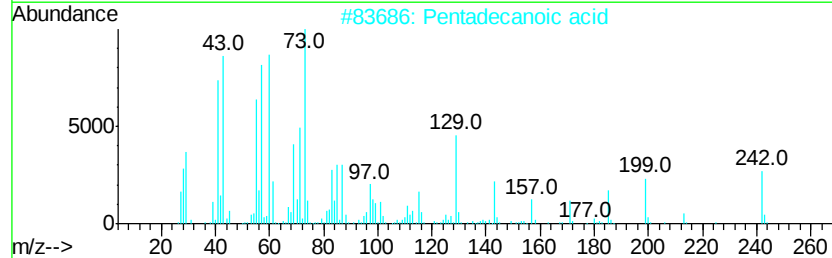
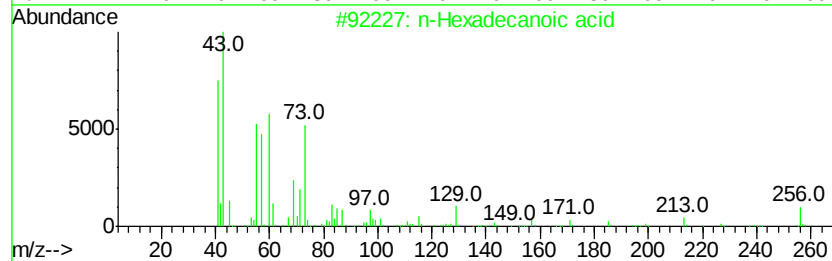
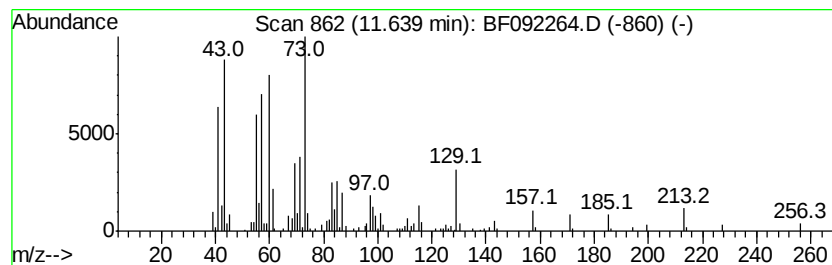
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	4.06 ng	314355	Phenanthrene-d10	11.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	83
3		Tridecanoic acid	214	C13H26O2	000638-53-9	80
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	59
5		n-Decanoic acid	172	C10H20O2	000334-48-5	59



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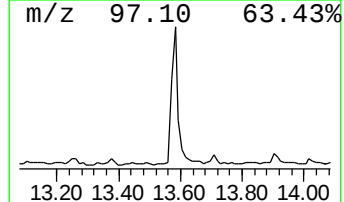
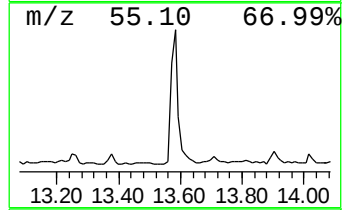
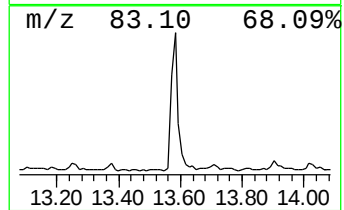
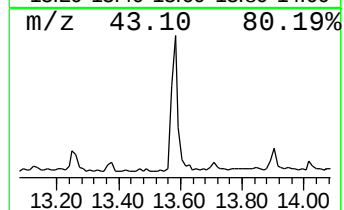
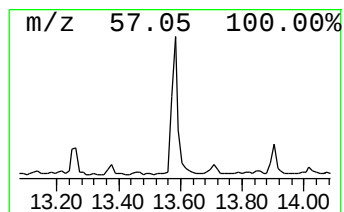
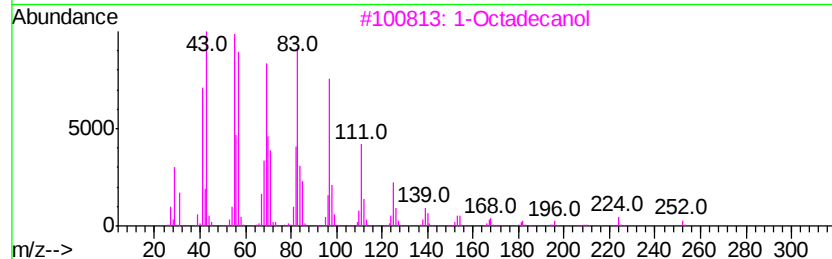
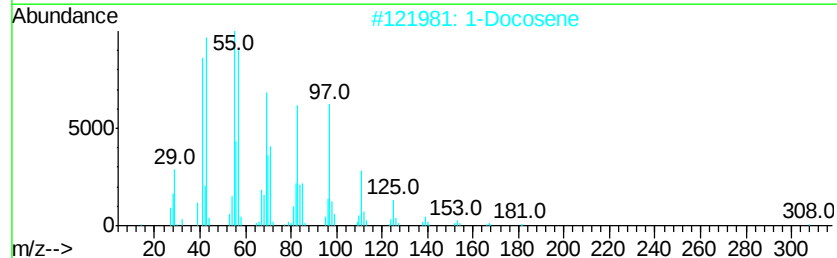
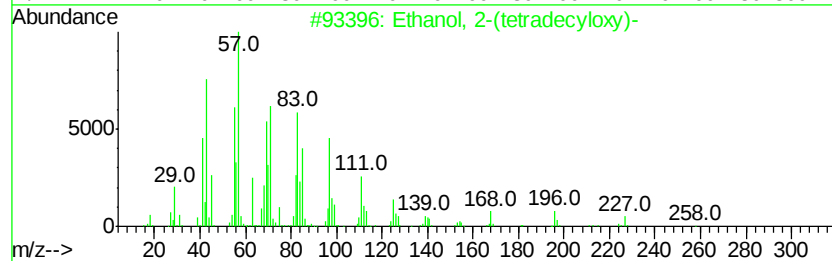
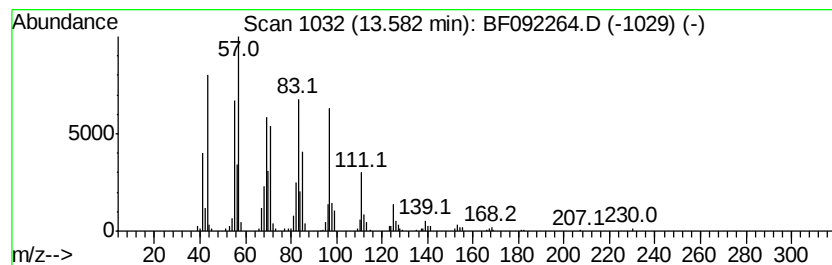
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Ethanol, 2-(tetradecyloxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.58	7.36 ng	473700	Chrysene-d12	13.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
2		1-Docosene	308	C22H44	001599-67-3	81
3		1-Octadecanol	270	C18H38O	000112-92-5	76
4		1-Nonadecene	266	C19H38	018435-45-5	76
5		1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	76



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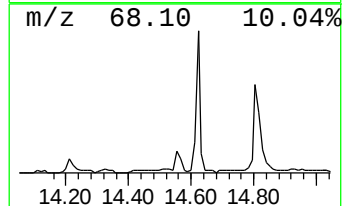
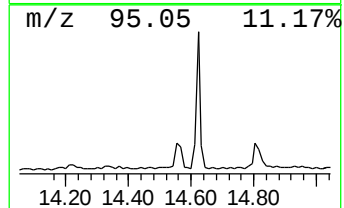
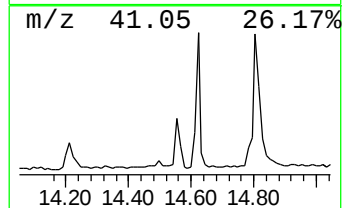
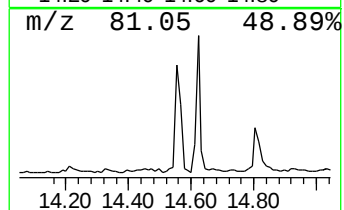
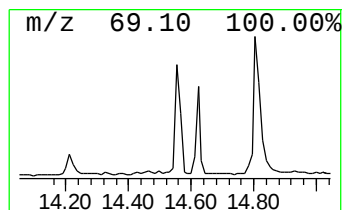
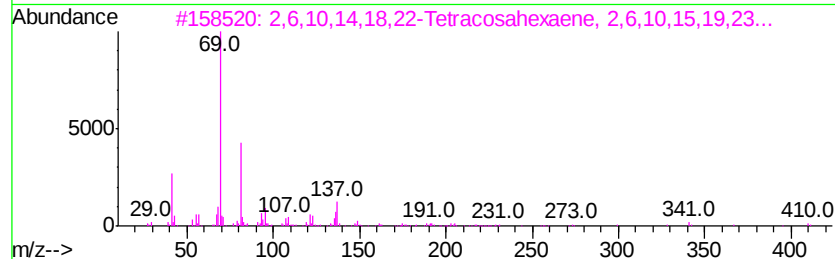
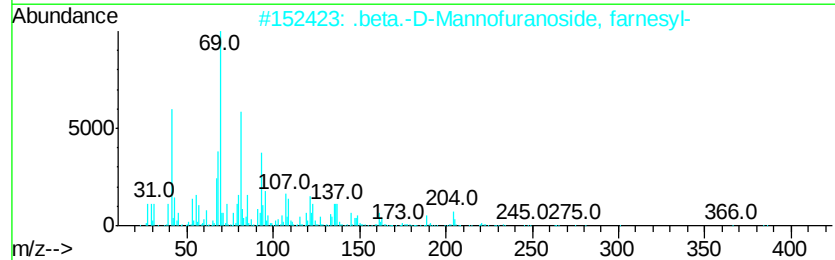
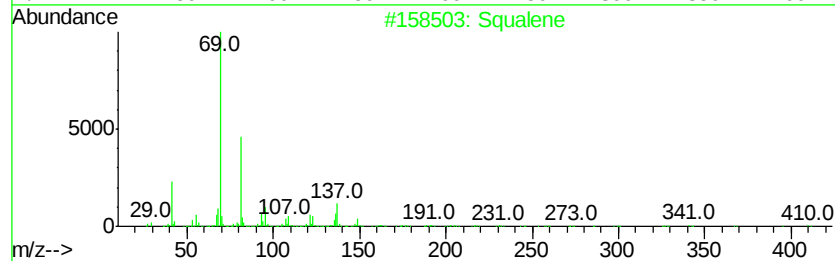
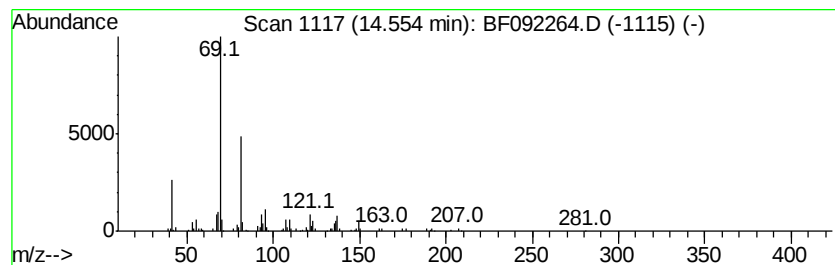
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TIC Integration Parameters: LSCINT.P

Peak Number 7 Squalene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.55	3.09 ng	140617	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	91
2		.beta.-D-Mannofuranoside, farnesyl-	384	C21H36O6	1000155-15-5	78
3		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	72
4		1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	64
5		2,6-Octadiene, 4-methyl-	124	C9H16	074498-94-5	58



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Data File : BF092264.D
Acq On : 14 Jan 2017 1:42
Operator : UM/SJ
Sample : I1031-13
Misc :
ALS Vial : 25 Sample Multiplier: 1

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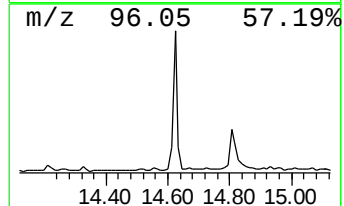
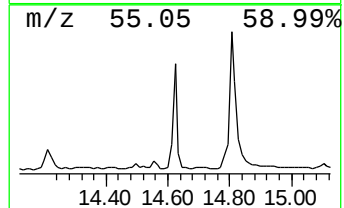
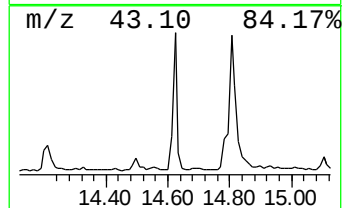
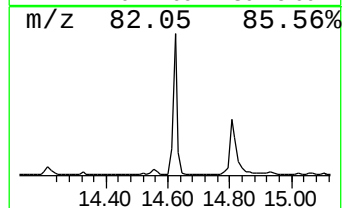
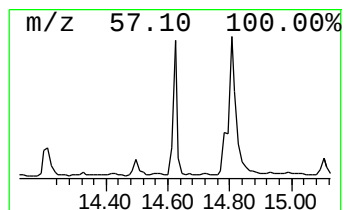
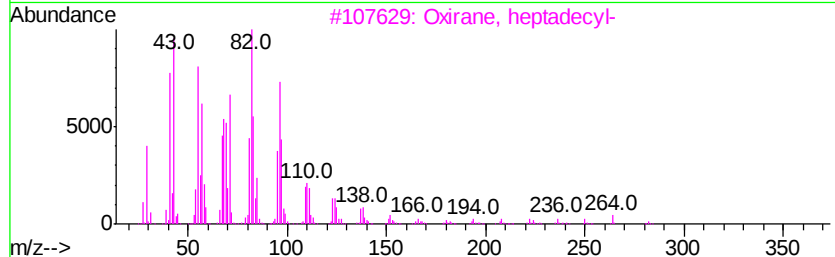
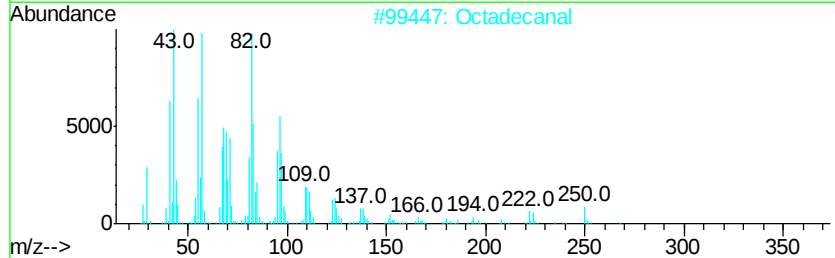
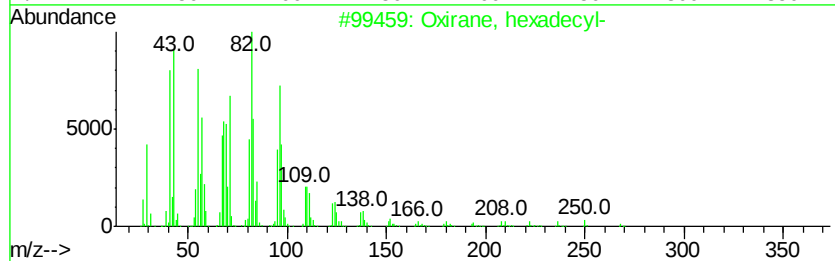
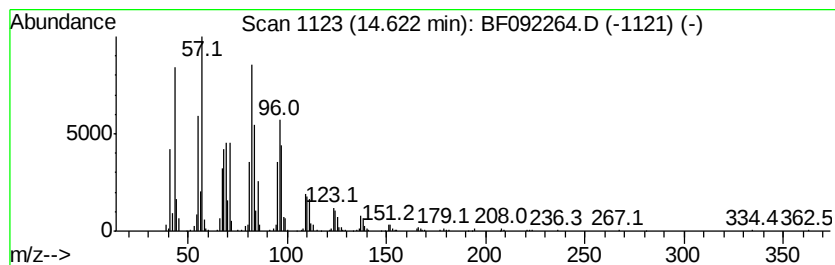
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Oxirane, hexadecyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.62	12.92 ng	588615	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
2		Octadecanal	268	C18H36O	000638-66-4	91
3		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
4		Tetracosanal	352	C24H48O	057866-08-7	91
5		17-Octadecenal	266	C18H34O	056554-86-0	90



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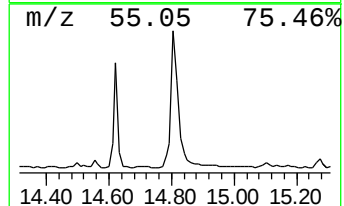
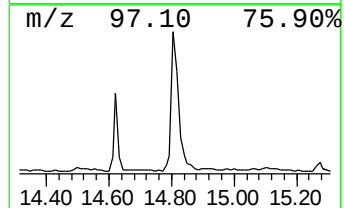
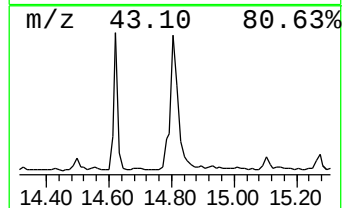
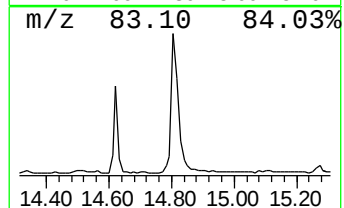
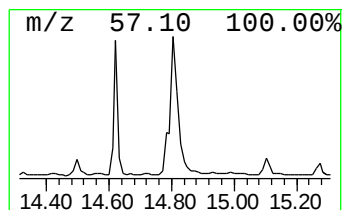
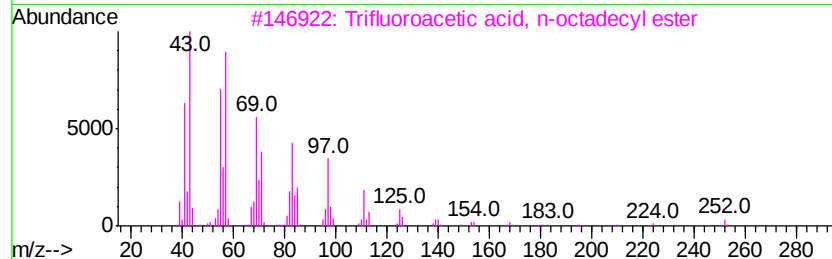
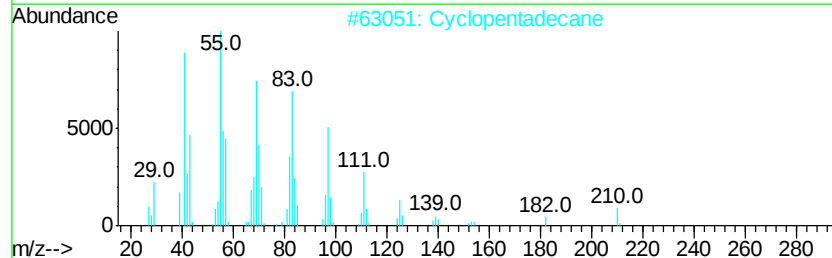
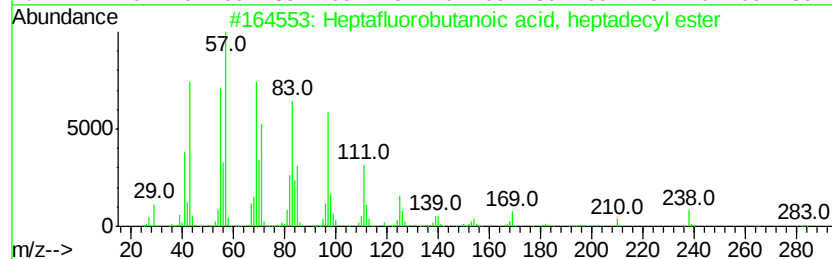
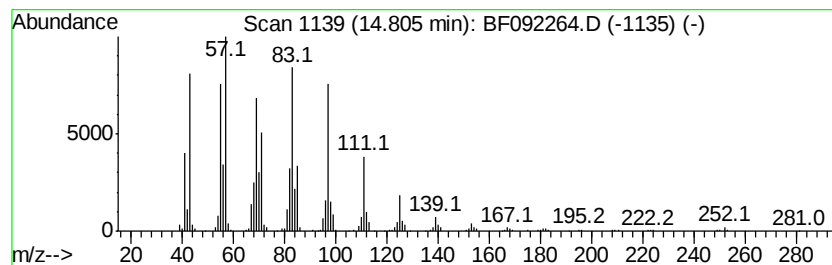
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Heptafluorobutanoic acid, h... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	22.98 ng	1046600	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
2		Cyclopentadecane	210	C15H30	000295-48-7	93
3		Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	91
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
5		1-Docosene	308	C22H44	001599-67-3	91



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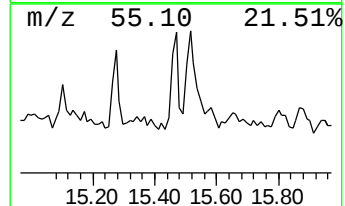
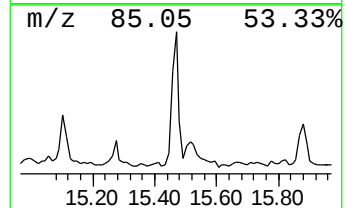
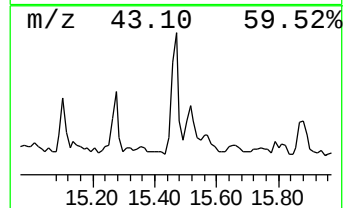
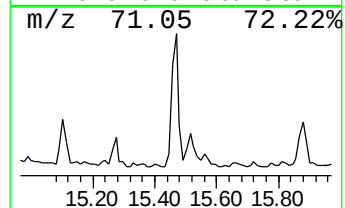
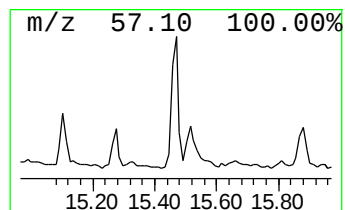
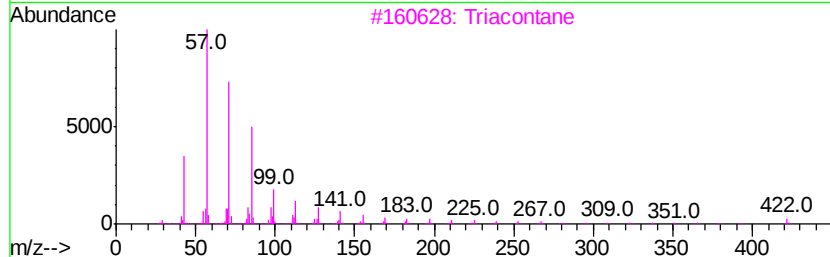
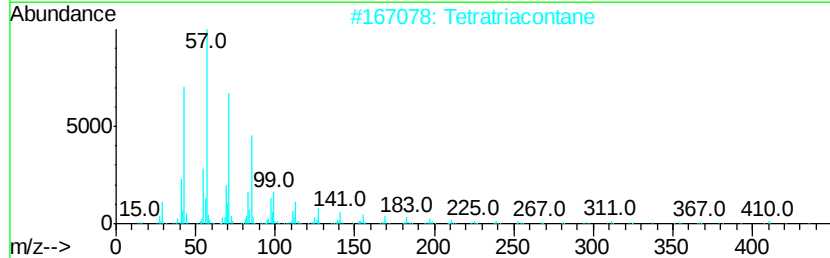
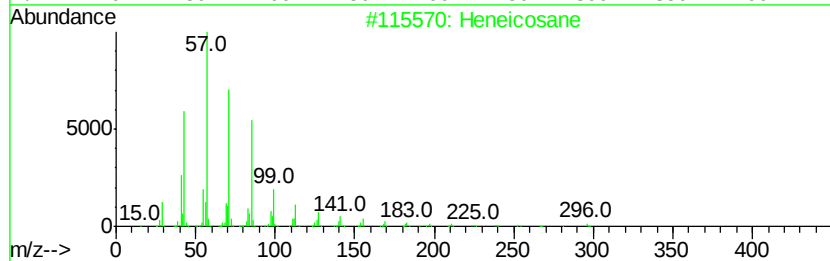
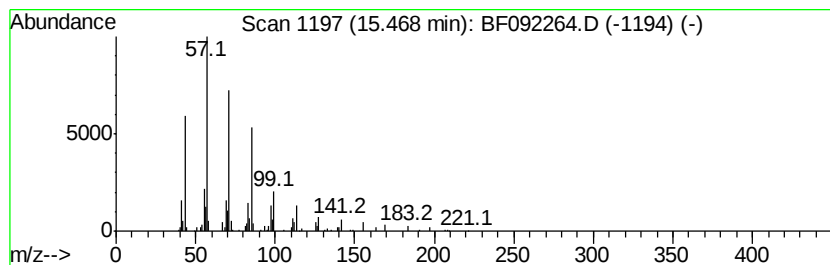
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.47	3.17 ng	144547	Perylene-d12	15.07	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Tetratriacontane	479	C34H70	014167-59-0	91
3	Triacontane	422	C30H62	000638-68-6	91
4	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
5	Tetratetracontane	619	C44H90	007098-22-8	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011317\
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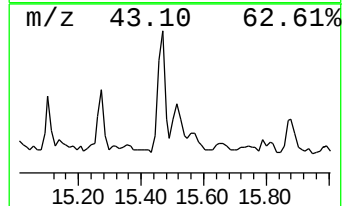
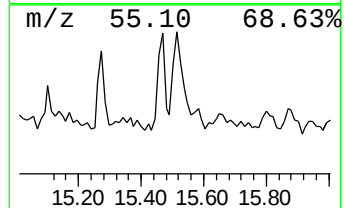
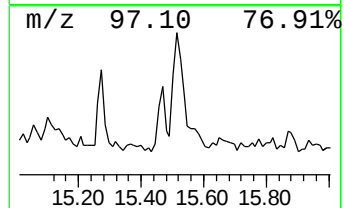
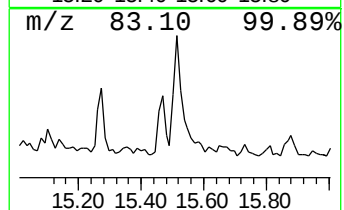
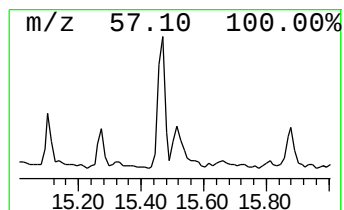
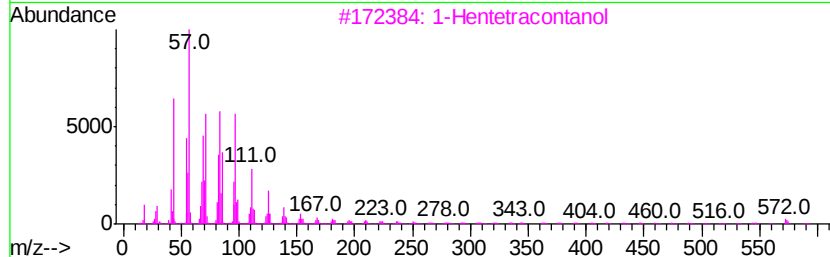
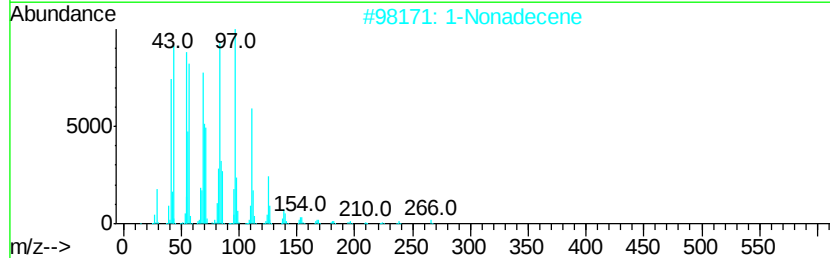
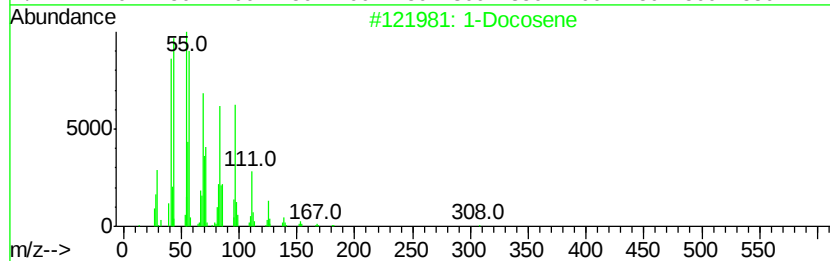
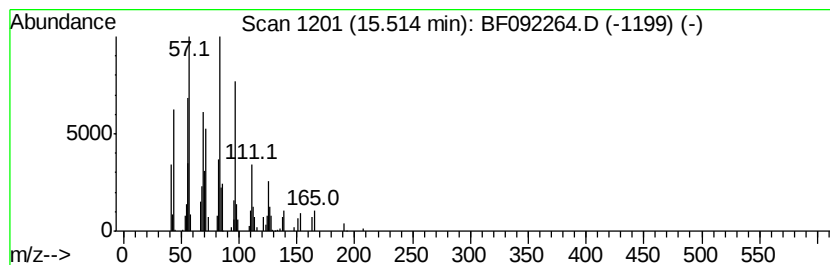
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Docosene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.51	2.18 ng	99293	Perylene-d12	15.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Docosene	308	C22H44	001599-67-3	64
2			1-Nonadecene	266	C19H38	018435-45-5	52
3			1-Hentetracontanol	593	C41H84O	040710-42-7	52
4			2- Bromopropionic acid, pentadec...	362	C18H35BrO2	1000292-44-2	46
5			Cyclohexanone, 3-ethyl-	126	C8H14O	022461-89-8	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011317\
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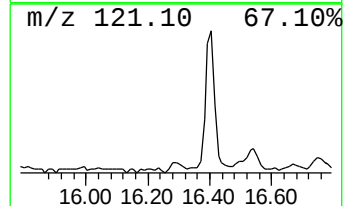
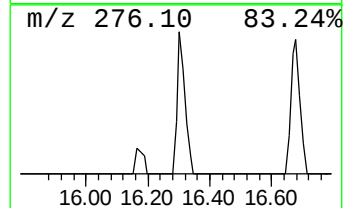
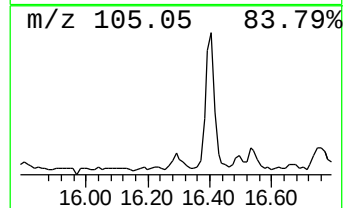
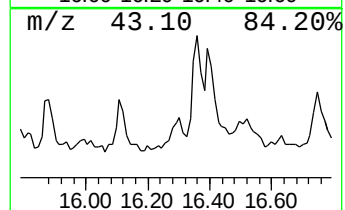
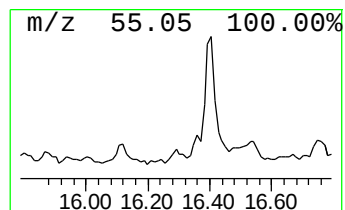
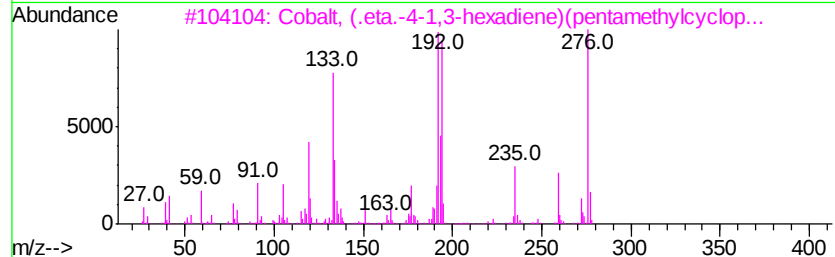
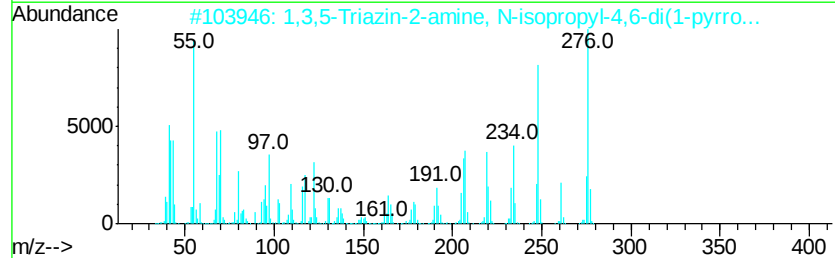
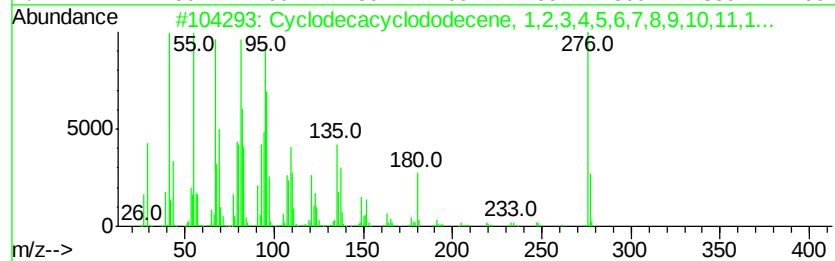
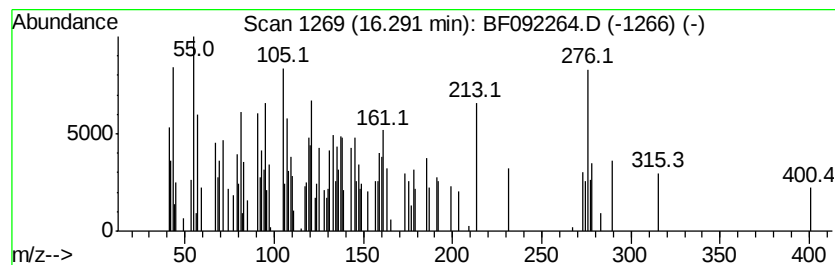
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown16.29 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.29	2.27 ng	103269	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclodecacyclododecene, 1,2,3,4,...	276	C20H36	014113-80-5	38
2		1,3,5-Triazin-2-amine, N-isoprop...	276	C14H24N6	1000276-80-0	35
3		Cobalt, (.eta.-4-1,3-hexadiene)(...	276	C16H25Co	1000156-29-9	22
4		Estran-3-one, 17-hydroxy-, (5.alpha....	276	C18H28O2	001434-85-1	14
5		Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011317\
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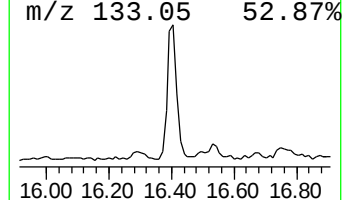
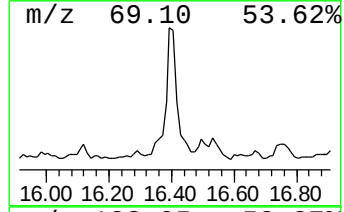
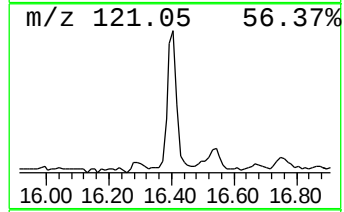
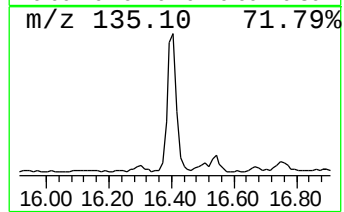
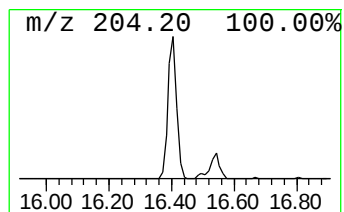
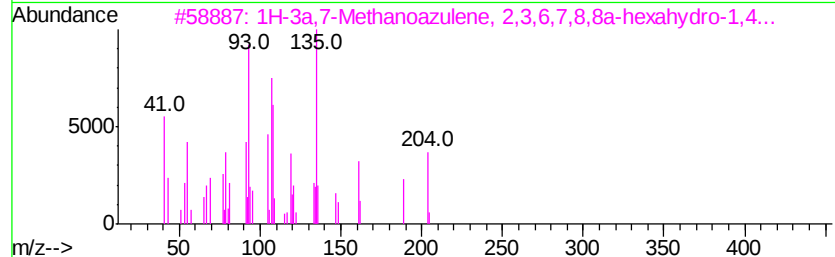
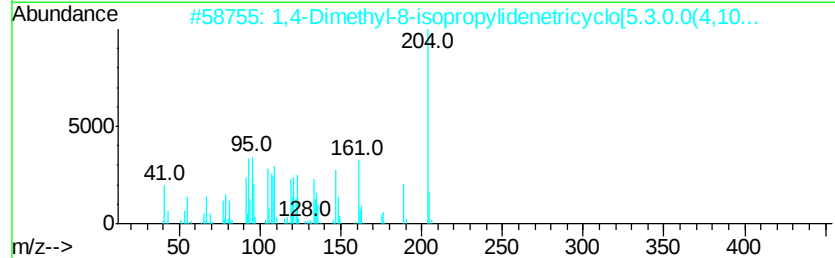
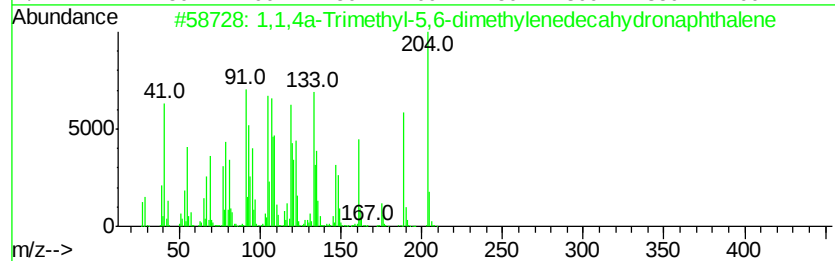
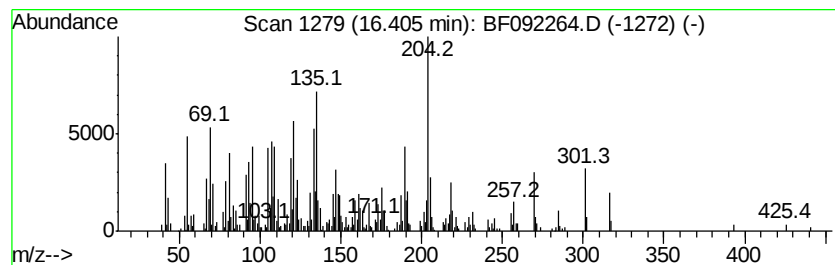
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 1,1,4a-Trimethyl-5,6-dimeth... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.41	25.89 ng	1179110	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	58
2		1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1000140-07-7	53
3		1H-3a,7-Methanoazulene, 2,3,6,7,...	204	C15H24	000560-32-7	53
4		Guaia-3,9-diene	204	C15H24	000489-83-8	47
5		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	004630-07-3	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011317\
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ClientSampleId :
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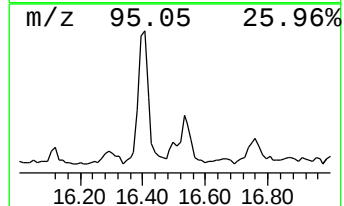
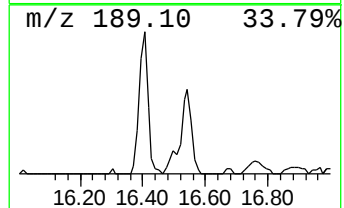
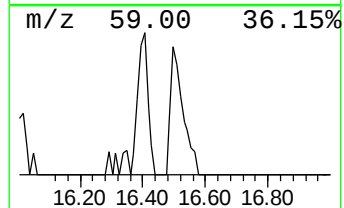
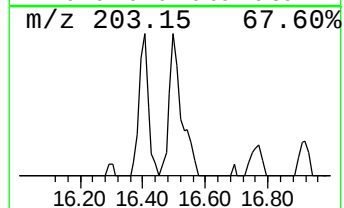
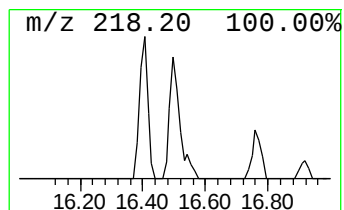
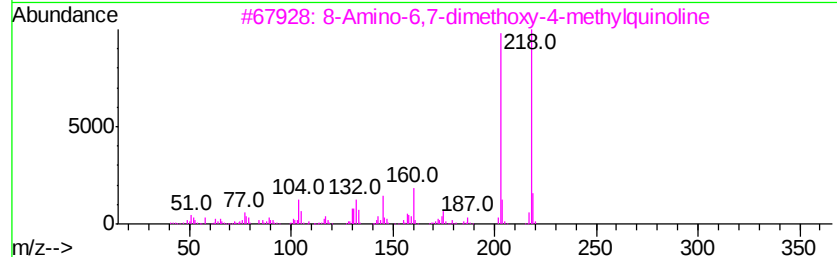
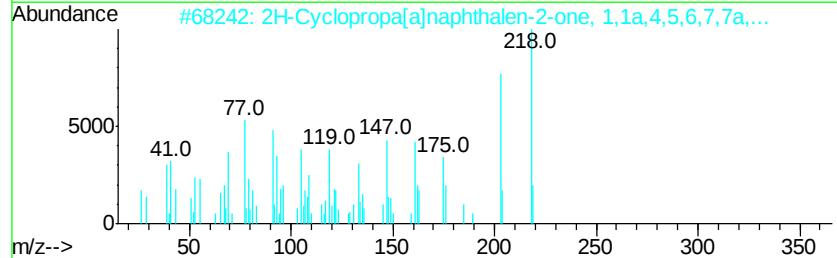
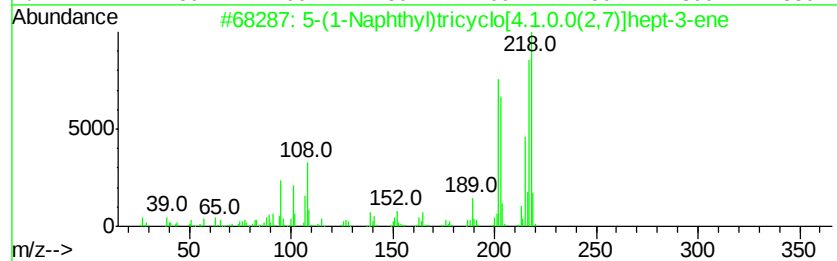
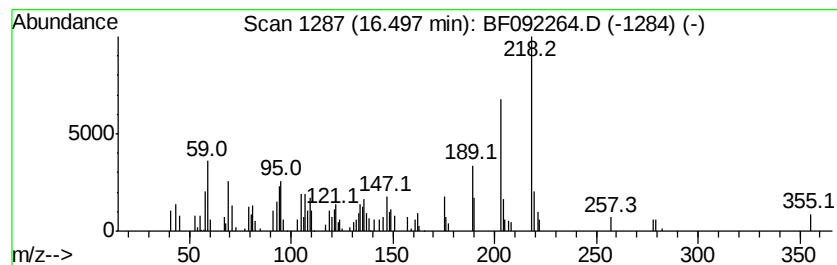
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 5-(1-Naphthyl)tricyclo[4.1.... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	2.42 ng	110376	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-(1-Naphthyl)tricyclo[4.1.0.0(2...]	218	C17H14	107729-05-5	58
2		2H-Cyclopropa[alnaphthalen-2-one...	218	C15H22O	006831-17-0	53
3		8-Amino-6,7-dimethoxy-4-methylqu...	218	C12H14N2O2	1000213-07-2	47
4		4,4,9,9-Tetramethyl-4,9-disilatr...	218	C12H18Si2	1000280-67-7	47
5		1-Hydroxypyrene	218	C16H10O	005315-79-7	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011317\
Data File : BF092264.D
Acq On : 14 Jan 2017 1:42
Operator : UM/SJ
Sample : I1031-13
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP-0027

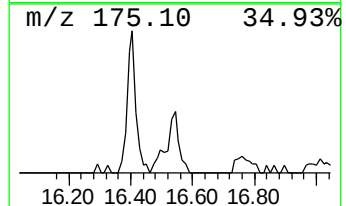
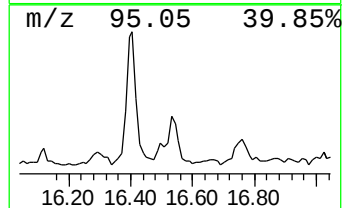
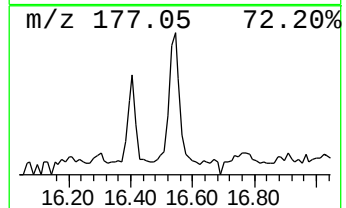
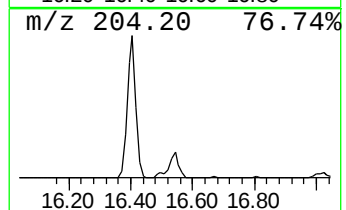
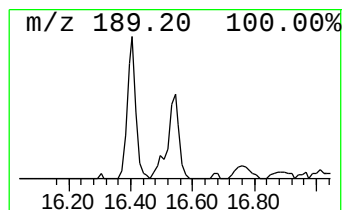
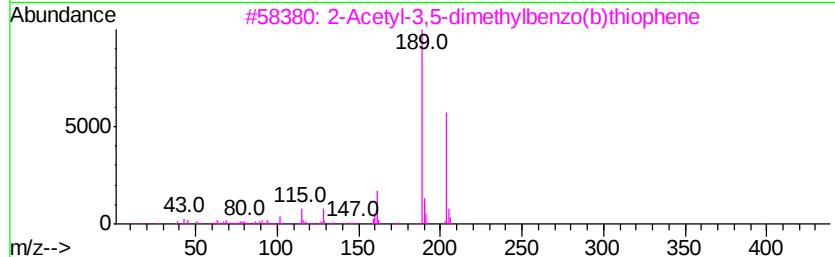
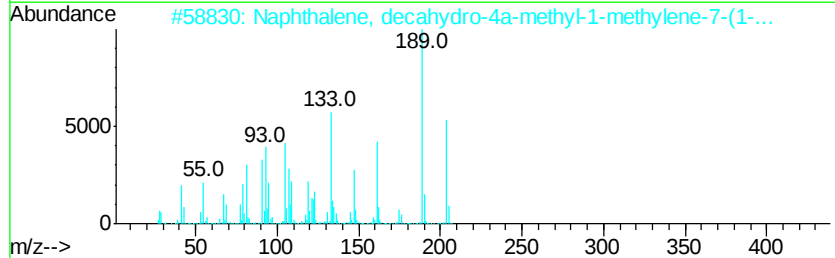
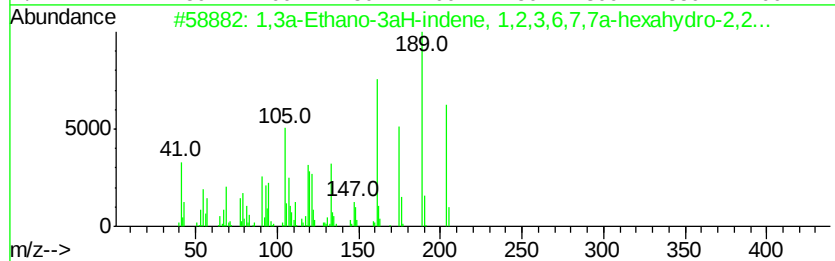
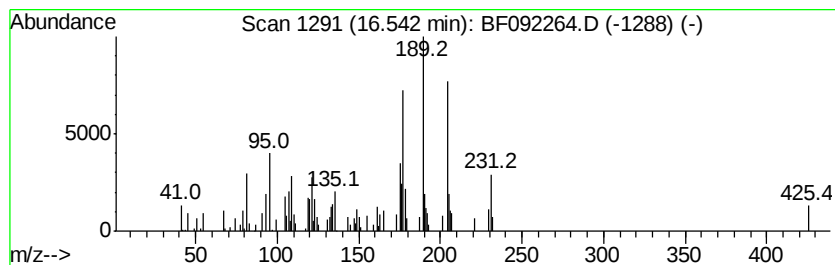
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown16.54 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.54	4.06 ng	185041	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3a-Ethano-3aH-indene, 1,2,3,6,...	204	C15H24	004545-68-0	49
2		Naphthalene, decahydro-4a-methyl...	204	C15H24	000515-17-3	47
3		2-Acetyl-3,5-dimethylbenzo(b)thi...	204	C12H12OS	006179-05-1	46
4		1H-Cyclopropa[a]naphthalene, 1a,...	204	C15H24	1000147-32-8	43
5		1H-Cyclopropa[a]naphthalene, 1a,...	204	C15H24	000489-29-2	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092264.D
Acq On : 14 Jan 2017 1:42
Operator : UM/SJ
Sample : I1031-13
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
COMP-0027

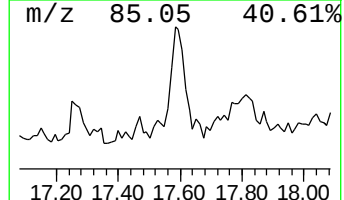
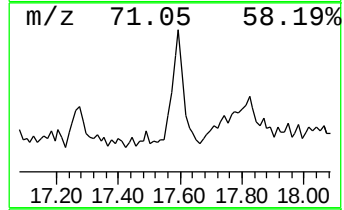
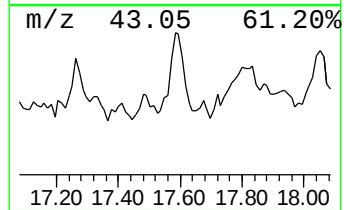
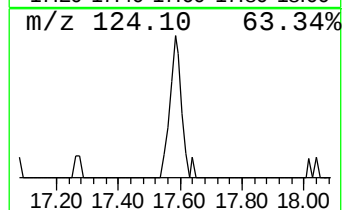
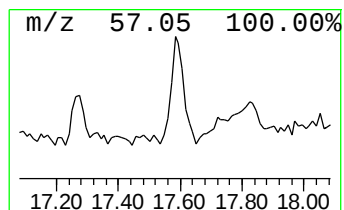
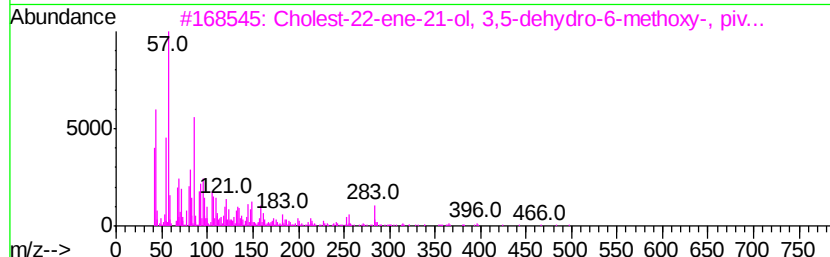
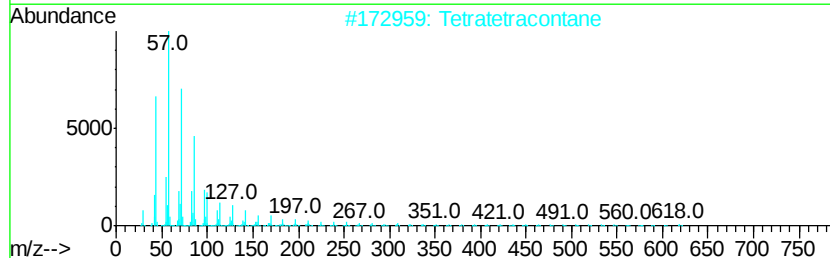
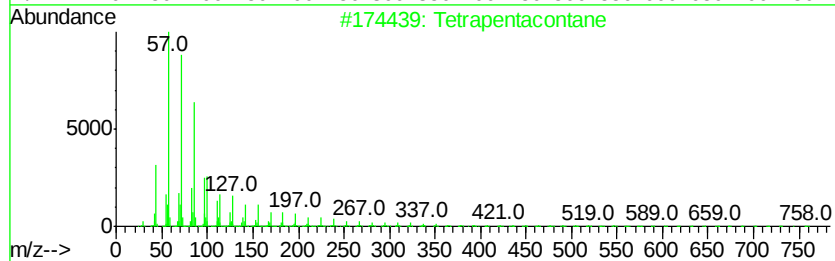
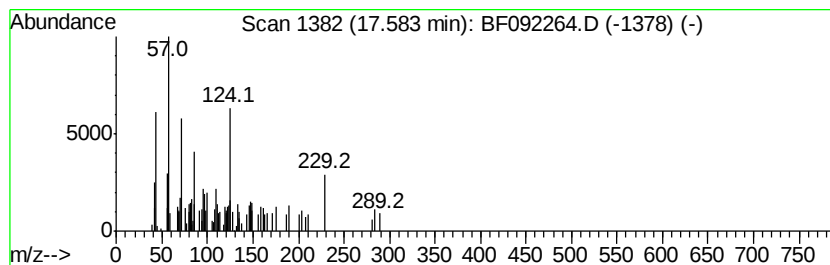
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown17.58 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.58	2.10 ng	95817	Perylene-d12	15.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrapentacontane	759	C54H110	005856-66-6	43
2			Tetratetracontane	619	C44H90	007098-22-8	43
3			Cholest-22-ene-21-ol, 3,5-dehydr...	498	C33H54O3	1000124-60-2	38
4			Pentacosane	352	C25H52	000629-99-2	30
5			Hexacosane	366	C26H54	000630-01-3	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092264.D
Acq On : 14 Jan 2017 1:42
Operator : UM/SJ
Sample : I1031-13
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
COMP-0027

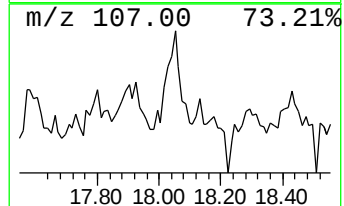
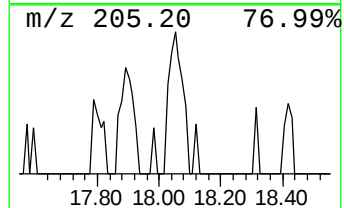
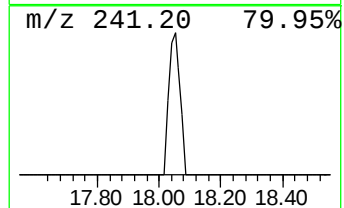
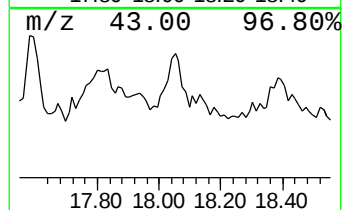
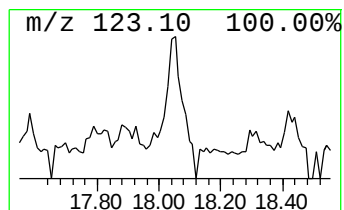
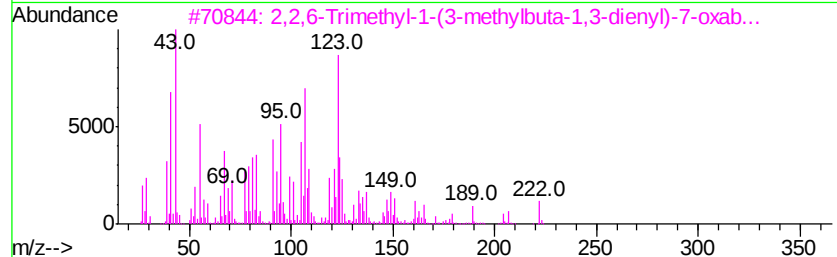
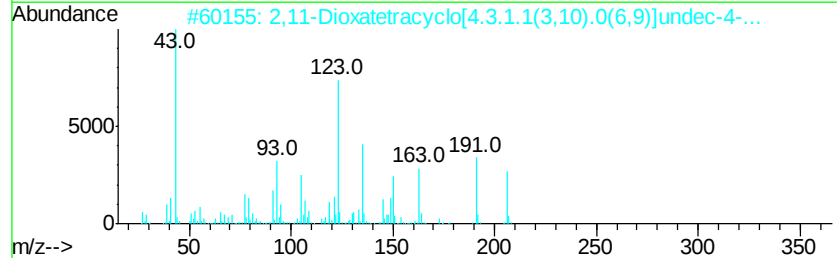
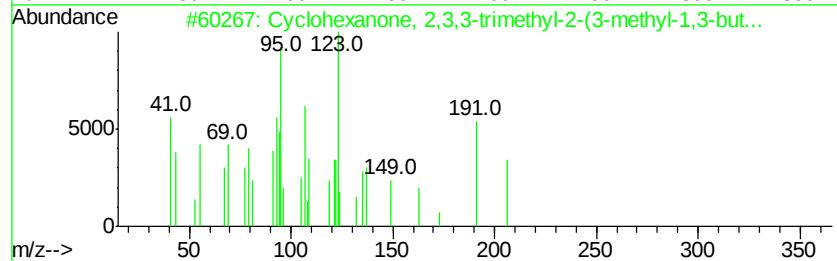
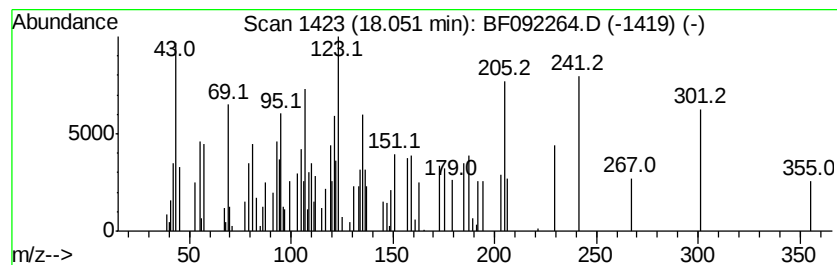
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Cyclohexanone, 2,3,3-trimet... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.05	2.02 ng	92182	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 2,3,3-trimethyl-2...	206	C14H22O	069296-90-8	53
2		2,11-Dioxatetracyclo[4.3.1.1(3,1...	206	C13H18O2	1000194-97-3	15
3		2,2,6-Trimethyl-1-(3-methylbuta...	222	C14H22O2	1000191-85-4	14
4		5-[1,2]Dithiolan-4-ylpentanamide	205	C8H15NOS2	1000195-38-7	10
5		Aniline, N-(3-butyl-2(Z),4-penta...	229	C16H23N	1000142-27-7	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011317\
 Data File : BF092264.D
 Acq On : 14 Jan 2017 1:42
 Operator : UM/SJ
 Sample : I1031-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-0027

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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...	4.70	39.7	ng	2334250	1	6.58	1176970	20.0
unknown6.35	6.35	84.3	ng	4962220	1	6.58	1176970	20.0
n-Hexadecanoic acid	11.64	4.1	ng	314355	4	11.08	1549910	20.0
Ethanol, 2-(tetra...	13.58	7.4	ng	473700	5	13.71	1286860	20.0
Squalene	14.55	3.1	ng	140617	6	15.07	910932	20.0
Oxirane, hexadecyl-	14.62	12.9	ng	588615	6	15.07	910932	20.0
Heptafluorobutano...	14.81	23.0	ng	1046600	6	15.07	910932	20.0
Heneicosane	15.47	3.2	ng	144547	6	15.07	910932	20.0
1-Docosene	15.51	2.2	ng	99293	6	15.07	910932	20.0
unknown16.29	16.29	2.3	ng	103269	6	15.07	910932	20.0
1,1,4a-Trimethyl-...	16.41	25.9	ng	1179110	6	15.07	910932	20.0
5-(1-Naphthyl)tri...	16.50	2.4	ng	110376	6	15.07	910932	20.0
unknown16.54	16.54	4.1	ng	185041	6	15.07	910932	20.0
2,2,6-Trimethyl-1...	16.76	4.0	ng	183685	6	15.07	910932	20.0
unknown17.58	17.58	2.1	ng	95817	6	15.07	910932	20.0
Cyclohexanone, 2,...	18.05	2.0	ng	92182	6	15.07	910932	20.0