

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040617\
 Data File : BF094261.D
 Acq On : 7 Apr 2017 8:44
 Operator : SJ/MA
 Sample : I2542-18
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-5

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-BF032217.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	3.993	353	358	366	rBV	626437	727742	19.85%	2.131%
2	4.393	421	426	429	rBV	3303133	3198767	87.24%	9.365%
3	4.781	488	492	498	rVB	43823	44261	1.21%	0.130%
4	5.063	537	540	543	rBV	77487	65713	1.79%	0.192%
5	5.463	602	608	613	rBV2	2912100	3580173	97.65%	10.482%
6	5.598	626	631	635	rBV	3322779	3455248	94.24%	10.116%
7	5.851	670	674	678	rBV	1069462	928736	25.33%	2.719%
8	5.945	686	690	695	rVB	99344	95791	2.61%	0.280%
9	6.034	700	705	708	rBV	2736826	2677085	73.02%	7.838%
10	6.504	779	785	788	rBV	1964761	2163011	58.99%	6.333%
11	7.045	873	877	888	rVB3	65859	110997	3.03%	0.325%
12	7.357	925	930	933	rBV	1221930	1247635	34.03%	3.653%
13	8.681	1149	1155	1158	rBV	3606225	3666460	100.00%	10.734%
14	9.175	1235	1239	1244	rBV	311000	276709	7.55%	0.810%
15	9.492	1286	1293	1299	rBV2	1208346	1321536	36.04%	3.869%
16	10.086	1390	1394	1399	rVB	61701	63134	1.72%	0.185%
17	10.469	1454	1459	1471	rBV	1801481	1940227	52.92%	5.680%
18	10.675	1490	1494	1502	rBV	82169	111222	3.03%	0.326%
19	11.227	1584	1588	1592	rBV	61410	59602	1.63%	0.174%
20	11.316	1599	1603	1607	rBV	1102260	1130186	30.82%	3.309%
21	12.427	1789	1792	1795	rBV3	46079	50224	1.37%	0.147%
22	12.810	1853	1857	1860	rVB	117926	117134	3.19%	0.343%
23	13.086	1899	1904	1908	rBV	92746	106593	2.91%	0.312%
24	13.133	1908	1912	1915	rBV2	58883	62171	1.70%	0.182%
25	13.298	1935	1940	1944	rBV	2306411	2411648	65.78%	7.061%
26	13.645	1997	1999	2002	rVB	42213	38287	1.04%	0.112%
27	14.010	2058	2061	2066	rBV5	30329	41849	1.14%	0.123%
28	14.068	2068	2071	2074	rVV	48188	51048	1.39%	0.149%
29	14.468	2134	2139	2145	rBV2	231916	413240	11.27%	1.210%
30	14.568	2151	2156	2160	rBV	832907	933324	25.46%	2.732%
31	14.868	2204	2207	2211	rBV	94522	95907	2.62%	0.281%
32	15.015	2229	2232	2236	rVB5	42199	44142	1.20%	0.129%
33	15.251	2268	2272	2280	rVB	259845	326615	8.91%	0.956%
34	15.398	2294	2297	2301	rBV	124300	127676	3.48%	0.374%

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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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35	15.610	2331	2333	2337	rVB2	94883	95806	2.61%	0.280%
36	15.762	2356	2359	2362	rBV2	73996	74539	2.03%	0.218%
37	15.798	2362	2365	2368	rBV	99949	122051	3.33%	0.357%
38	15.968	2390	2394	2404	rBV	303061	463773	12.65%	1.358%
39	16.198	2429	2433	2441	rVB	619933	757167	20.65%	2.217%
40	16.321	2451	2454	2457	rVB	71448	78112	2.13%	0.229%
41	16.498	2481	2484	2488	rVB3	105661	131001	3.57%	0.384%
42	16.721	2518	2522	2526	rBV	109653	148529	4.05%	0.435%
43	16.768	2526	2530	2535	rVV	157632	246102	6.71%	0.721%
44	17.427	2638	2642	2653	rVB4	85875	155624	4.24%	0.456%
45	18.086	2750	2754	2764	rBV8	83492	199999	5.45%	0.586%

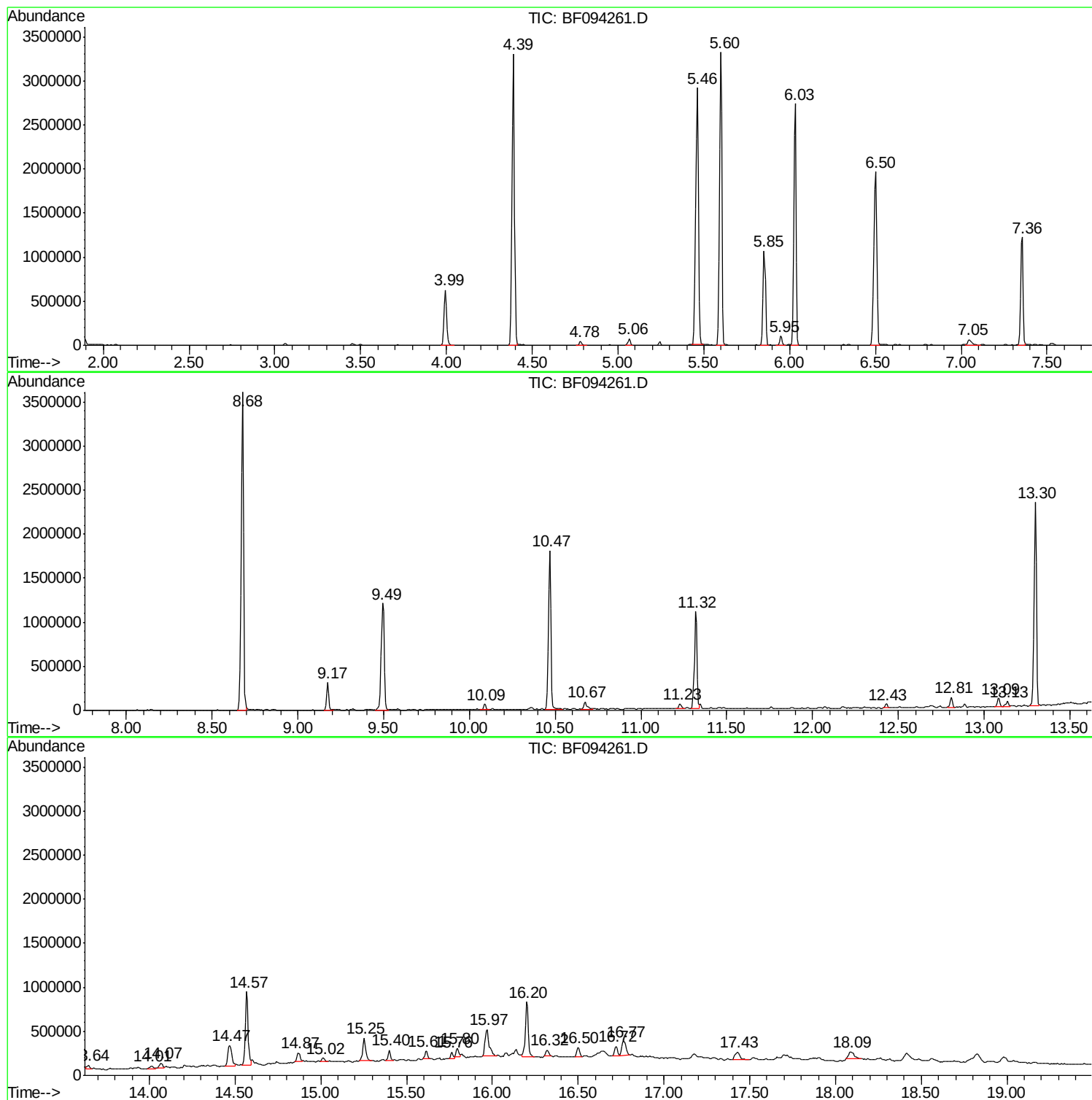
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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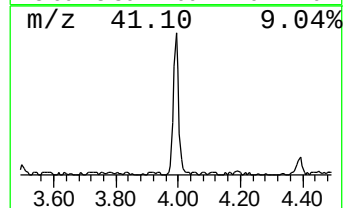
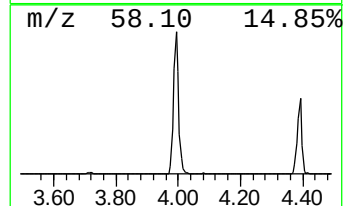
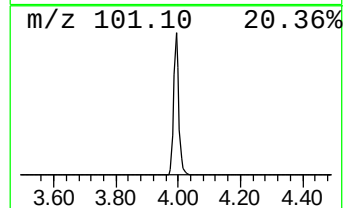
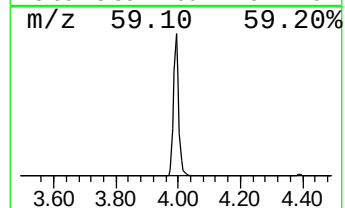
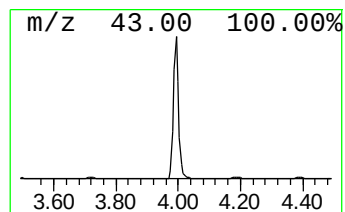
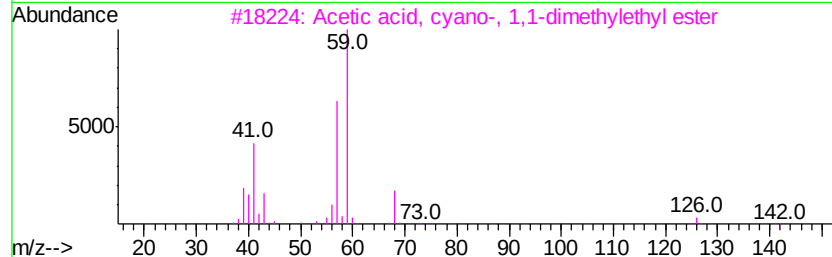
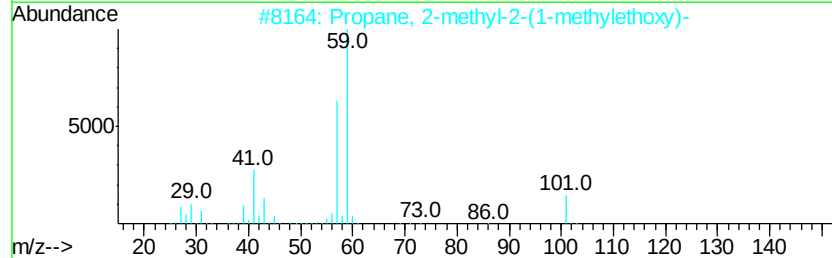
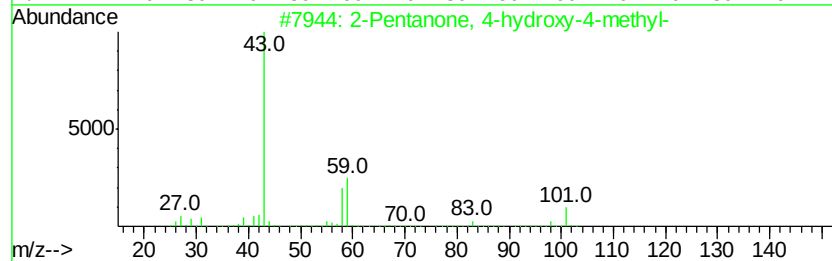
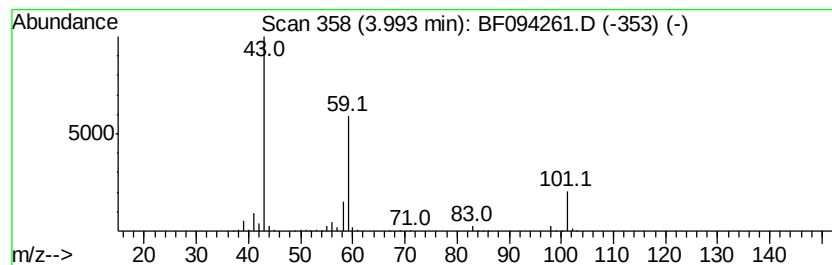
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
3.99	15.67 ng	727742	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	42
3		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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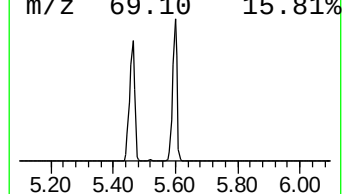
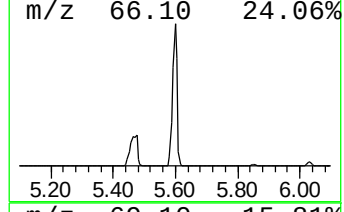
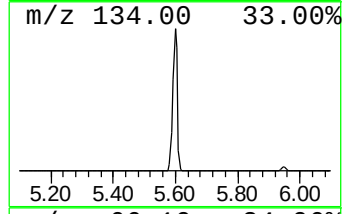
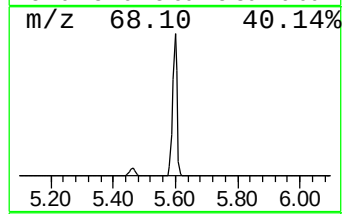
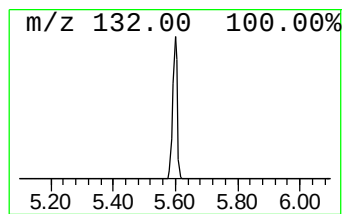
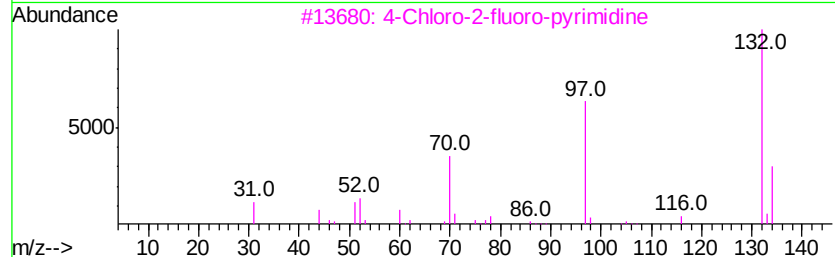
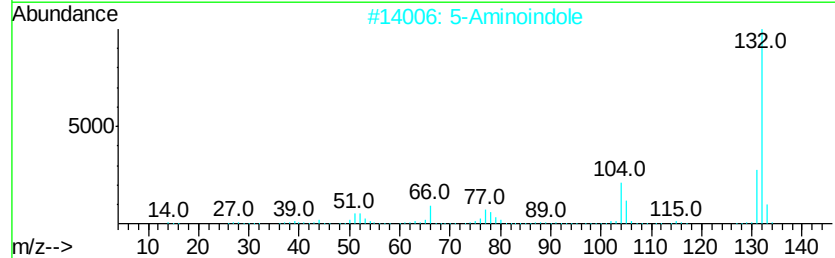
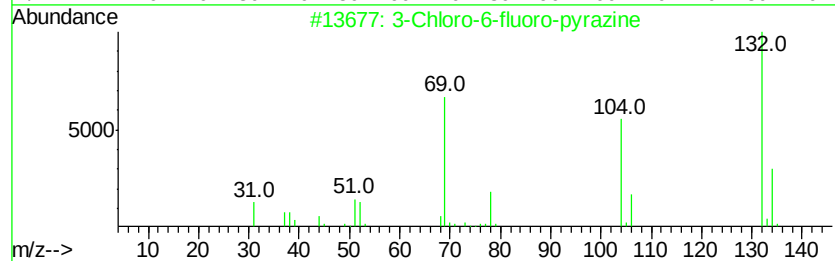
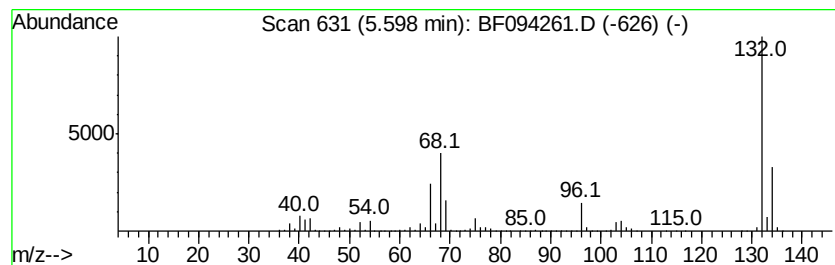
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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown5.60 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.60	74.41 ng	3455250	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5	Aminoindole	132	C8H8N2	005192-03-0	12
3	4	Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4	1	Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5	Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9	9



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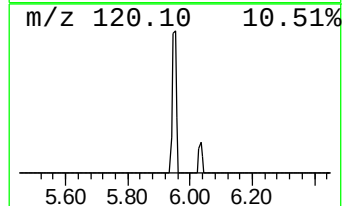
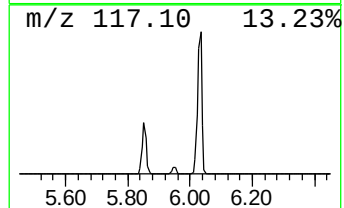
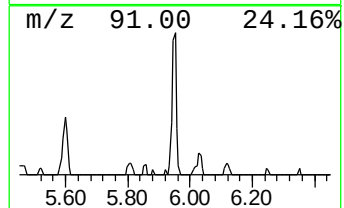
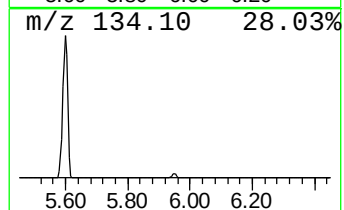
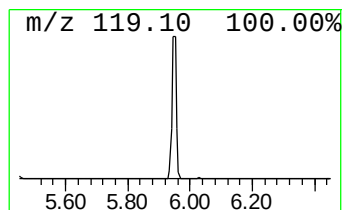
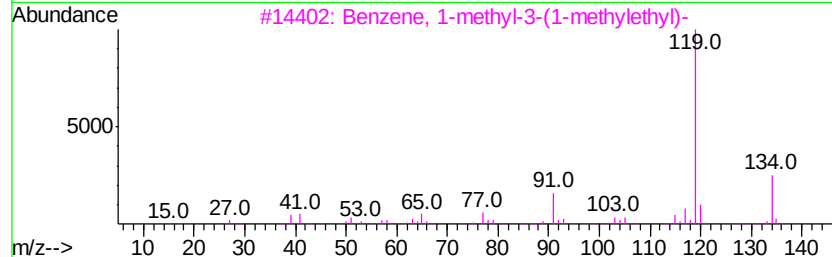
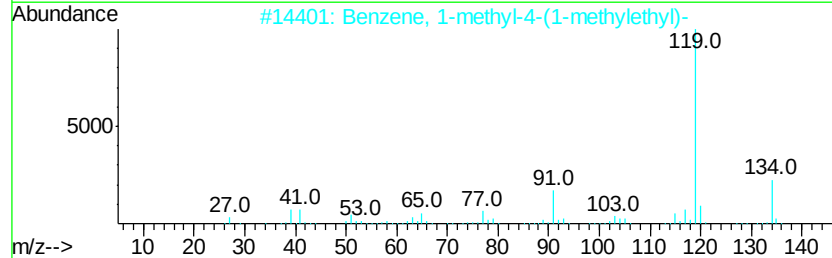
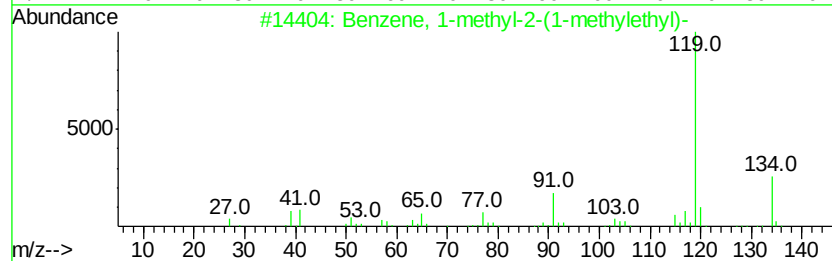
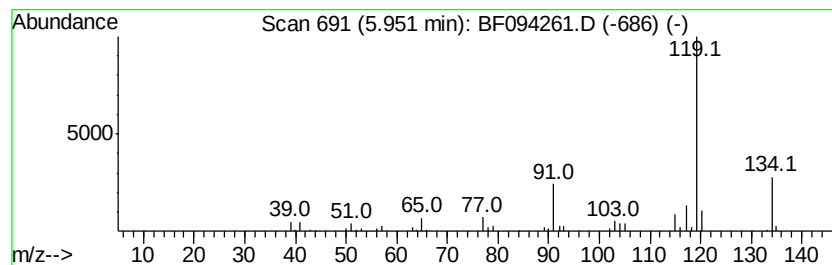
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Benzene, 1-methyl-2-(1-meth... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.95	2.06 ng	95791	1,4-Dichlorobenzene-d4	5.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	97
2		Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	97
3		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



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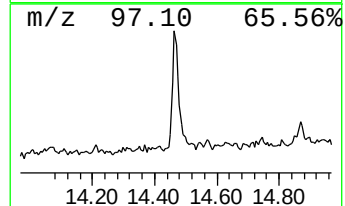
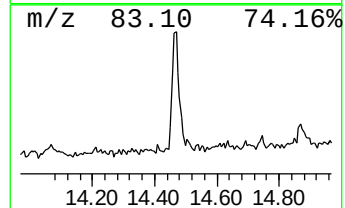
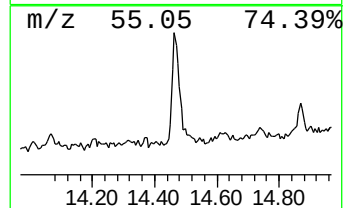
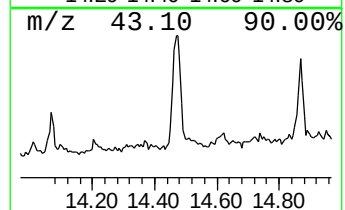
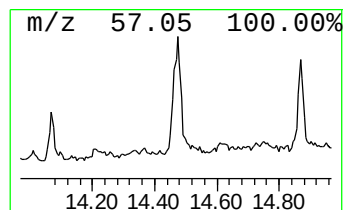
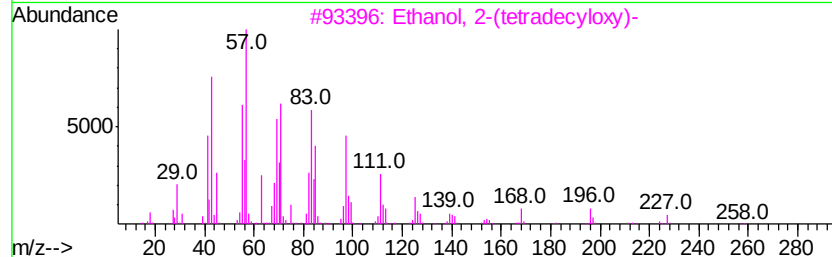
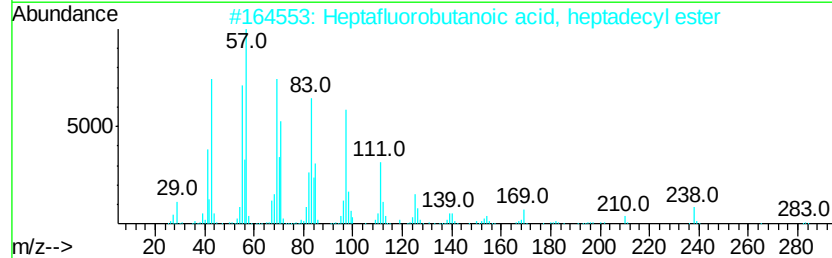
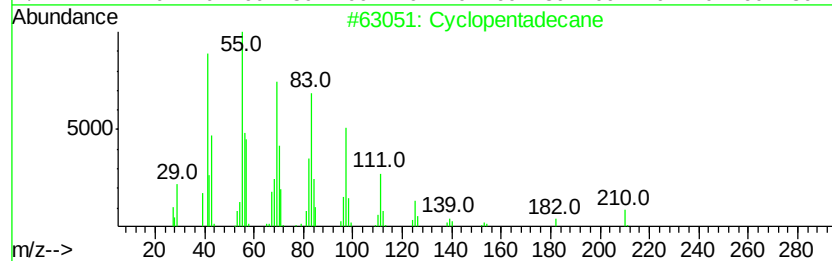
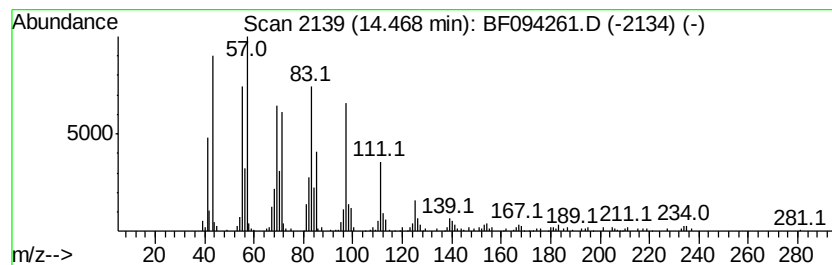
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Cyclopentadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.47	8.86 ng	413240	Chrysene-d12	14.57	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	93
2	Heptafluorobutanoic acid, heptadecanoic acid, heptadecanoic acid	452	C21H35F7O2	1000282-97-3	93
3	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
4	17-Pentatriacontene	491	C35H70	006971-40-0	91
5	Pentafluoropropionic acid, heptadecanoic acid, heptadecanoic acid	402	C20H35F5O2	1000283-04-2	90



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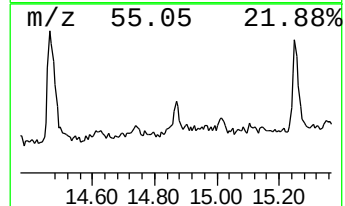
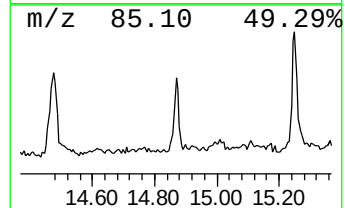
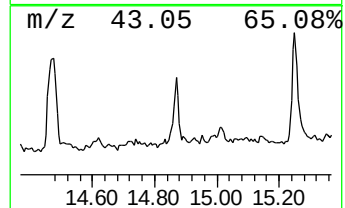
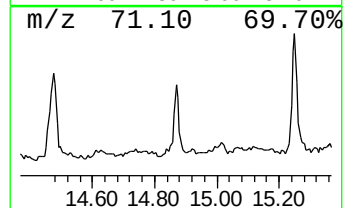
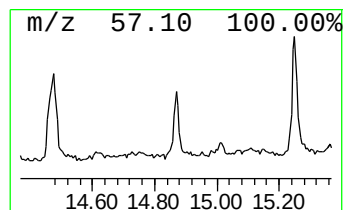
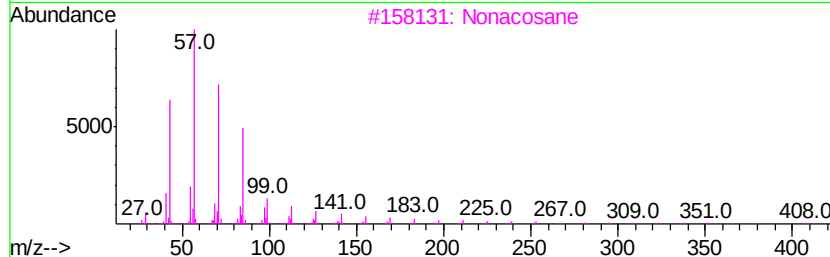
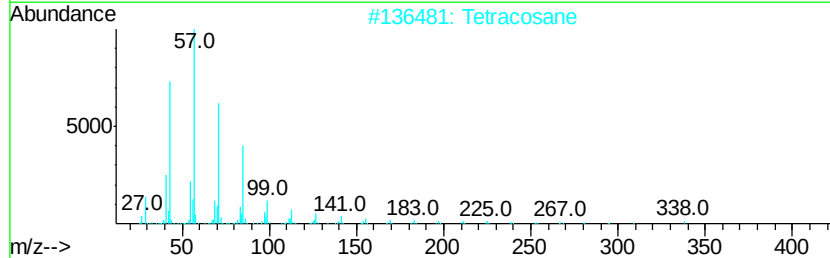
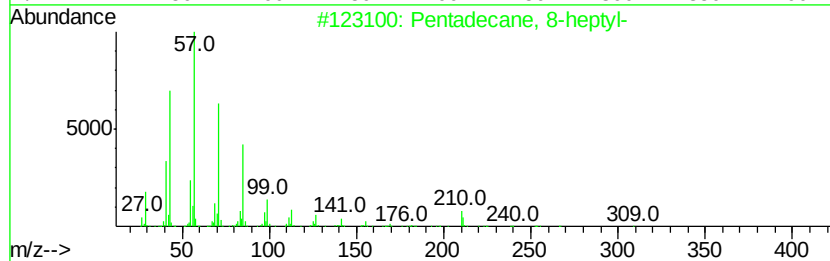
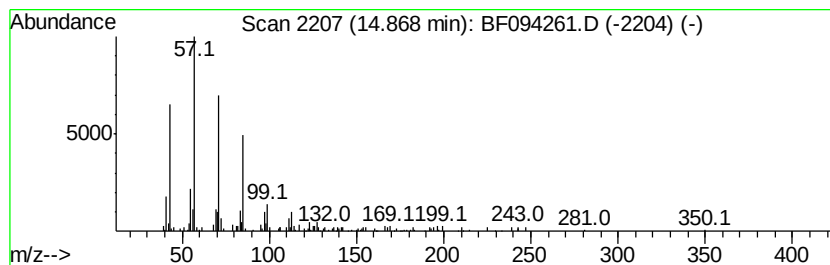
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Pentadecane, 8-heptyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	2.06 ng	95907	Chrysene-d12	14.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	86
2		Tetracosane	338	C24H50	000646-31-1	86
3		Nonacosane	408	C29H60	000630-03-5	86
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
5		Benzeneacetic acid, 2-tetradecyl...	332	C22H36O2	1000282-02-4	78



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Operator : SJ/MA
Sample : I2542-18
Misc :
ALS Vial : 43 Sample Multiplier: 1

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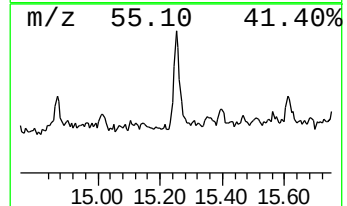
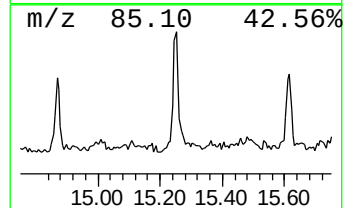
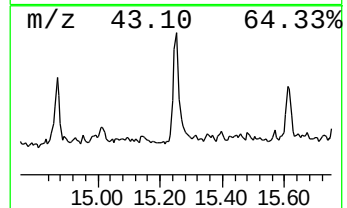
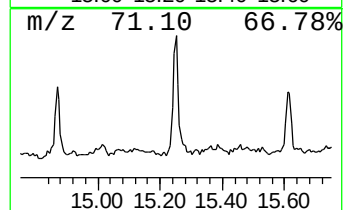
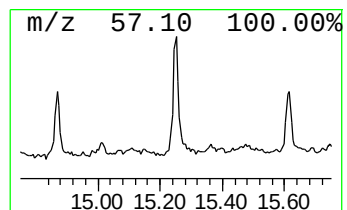
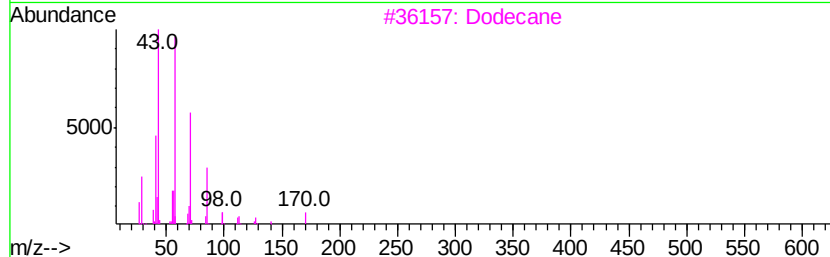
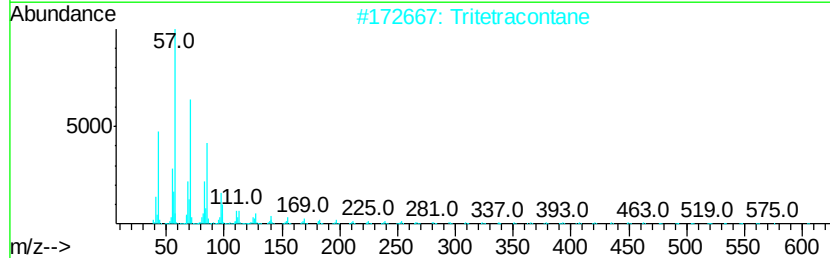
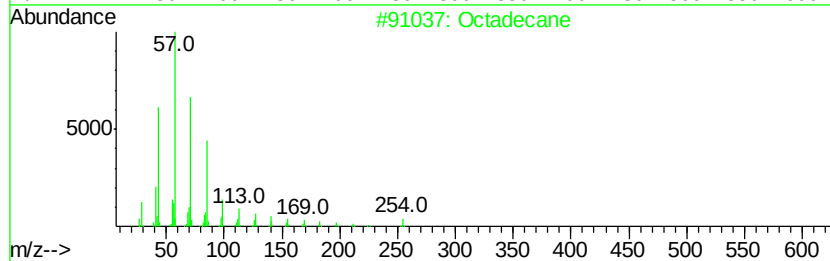
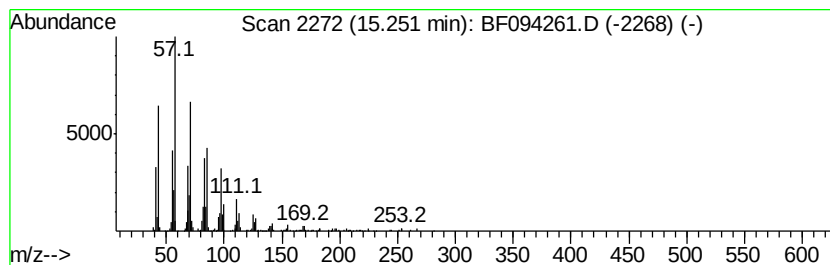
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.25	7.00 ng	326615	Chrysene-d12	14.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane		254	C18H38	000593-45-3	86
2	Tritetracontane		605	C43H88	007098-21-7	83
3	Dodecane		170	C12H26	000112-40-3	64
4	Heptafluorobutyric acid, n-octad...		466	C22H37F7O2	000400-57-7	60
5	Hexadecane		226	C16H34	000544-76-3	58



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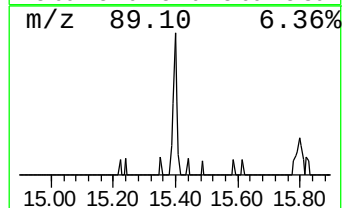
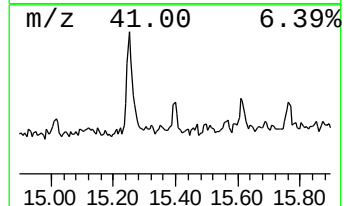
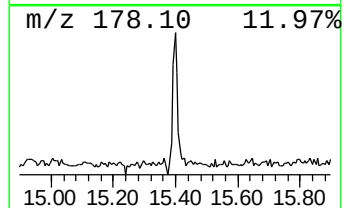
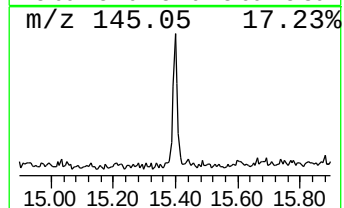
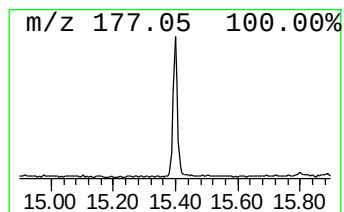
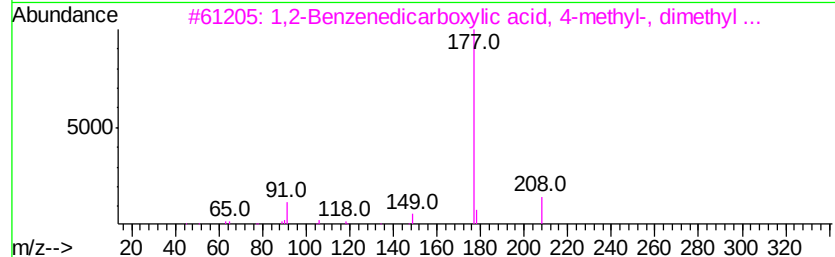
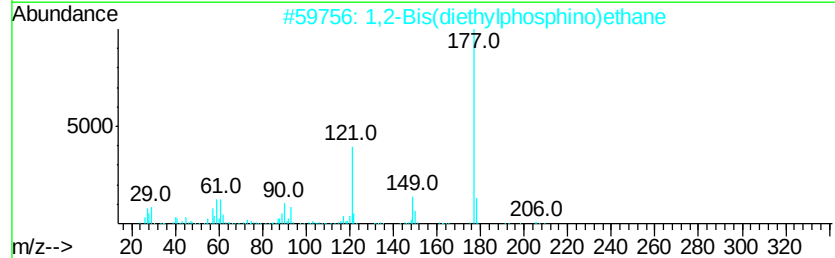
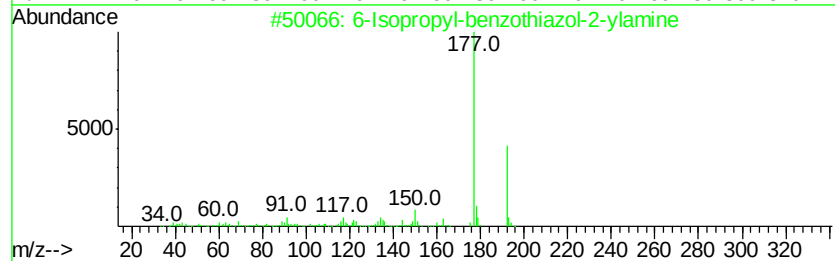
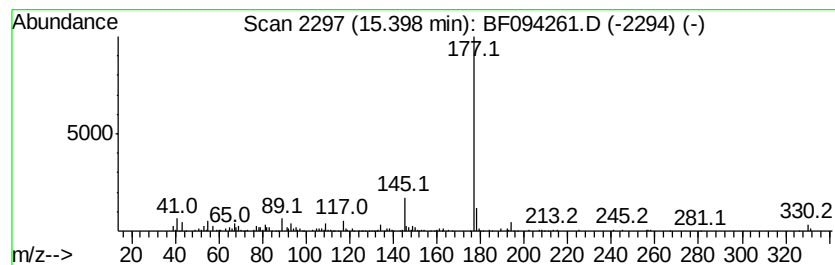
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown15.40 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.40	3.37 ng	127676	Perylene-d12	16.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Isopropyl-benzothiazol-2-ylamine	192	C10H12N2S	032895-14-0	38
2			1,2-Bis(diethylphosphino)ethane	206	C10H24P2	006411-21-8	38
3			1,2-Benzenedicarboxylic acid, 4-...	208	C11H12O4	020116-65-8	36
4			2H-1-Benzopyran, 3,5,6,8a-tetra...	192	C13H20O	041678-29-9	36
5			Phenol, 2,6-bis(1-methylpropyl)-	206	C14H22O	005510-99-6	36



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040617\
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Sample : I2542-18
Misc :
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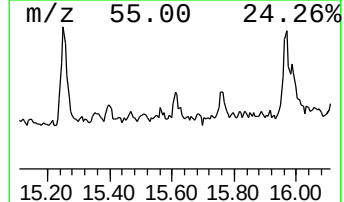
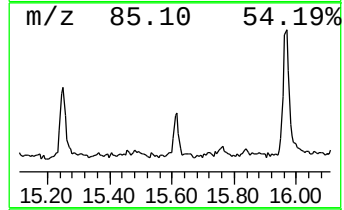
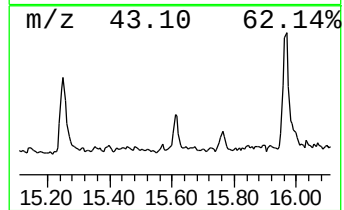
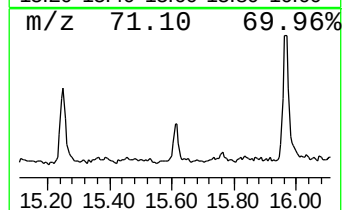
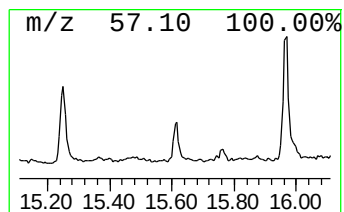
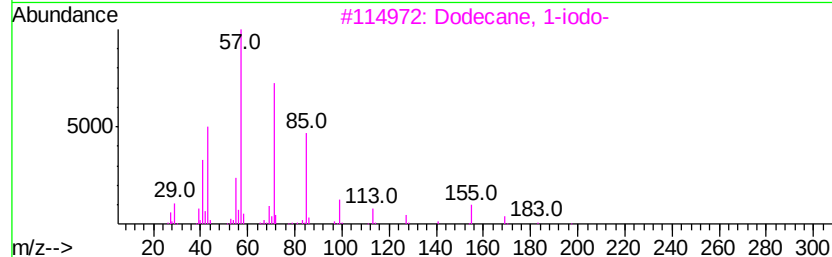
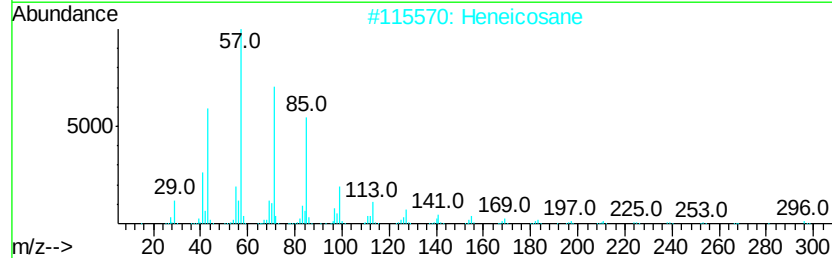
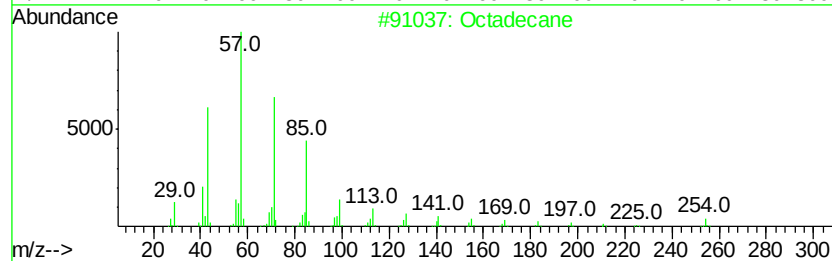
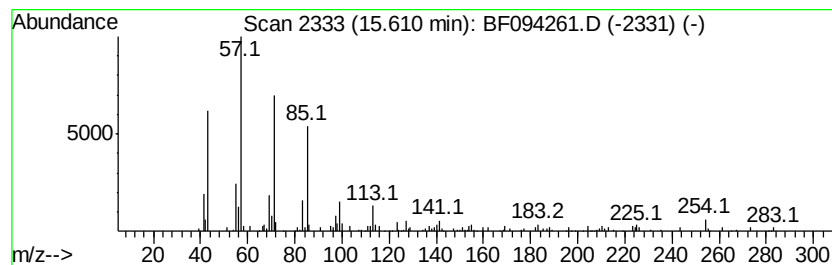
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Octadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	2.53 ng	95806	Perylene-d12	16.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	93
2		Heneicosane	296	C21H44	000629-94-7	83
3		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
4		Docosane	310	C22H46	000629-97-0	80
5		Eicosane	282	C20H42	000112-95-8	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040617\
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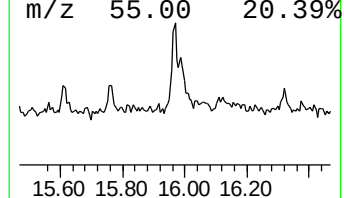
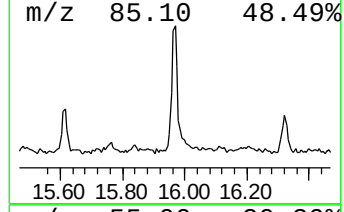
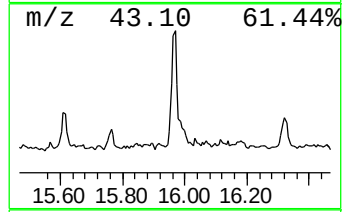
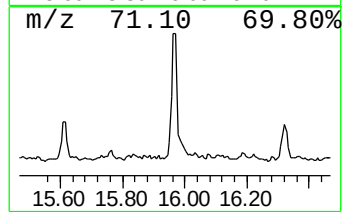
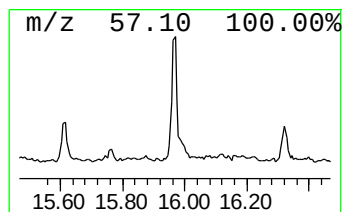
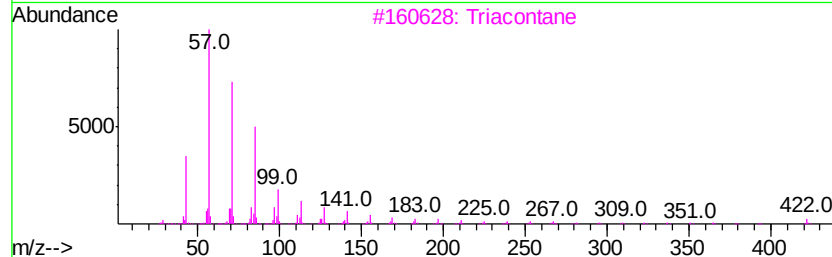
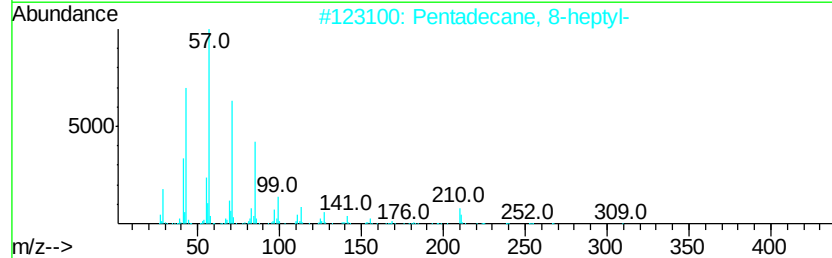
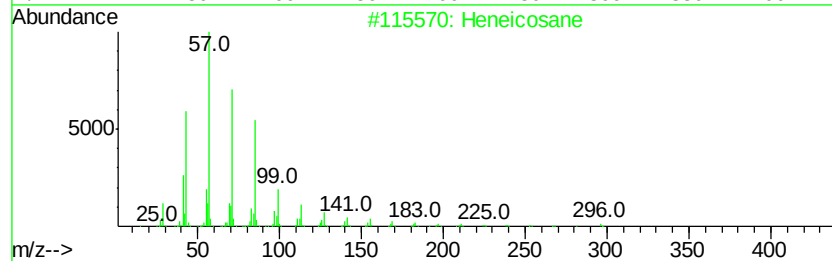
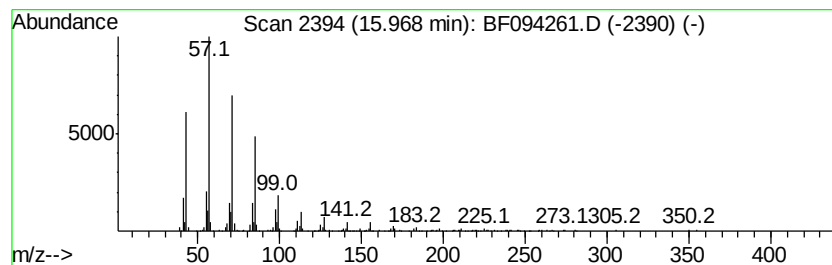
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	12.25 ng	463773	Perylene-d12	16.20

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Pentadecane, 8-heptyl-		310	C22H46	071005-15-7	91
3	Triacontane		422	C30H62	000638-68-6	90
4	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	86
5	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040617\
Data File : BF094261.D
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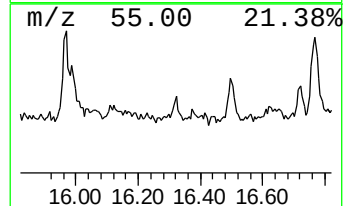
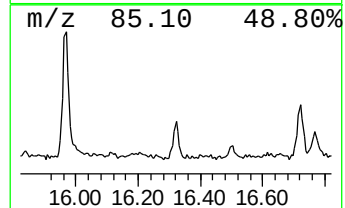
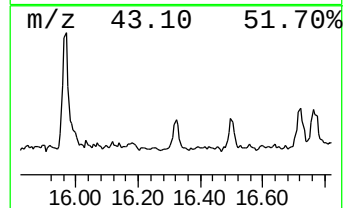
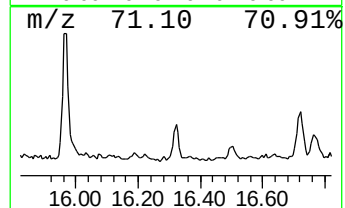
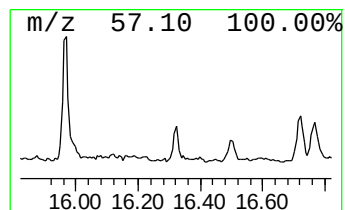
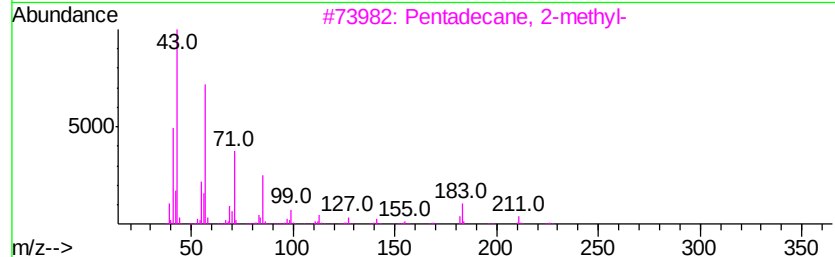
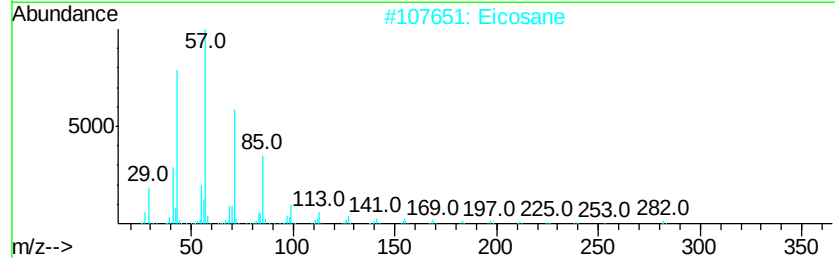
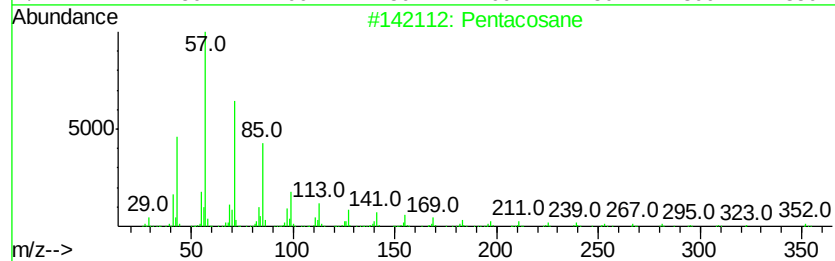
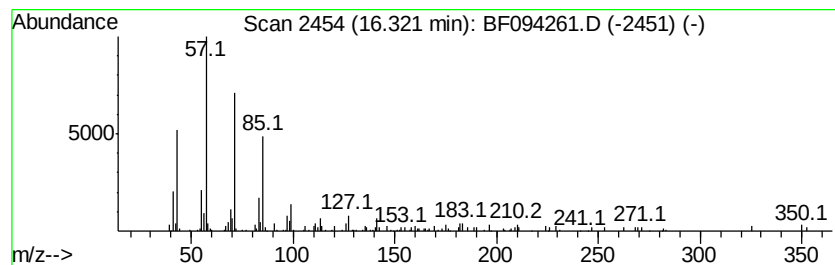
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Pentacosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.32	2.06 ng	78112	Perylene-d12	16.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacosane	352	C25H52	000629-99-2	93
2		Eicosane	282	C20H42	000112-95-8	90
3		Pentadecane, 2-methyl-	226	C16H34	001560-93-6	87
4		Nonadecane	268	C19H40	000629-92-5	86
5		Hexadecane	226	C16H34	000544-76-3	83



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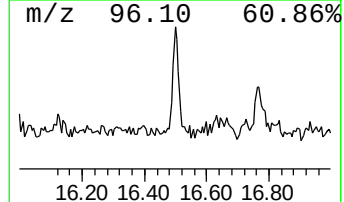
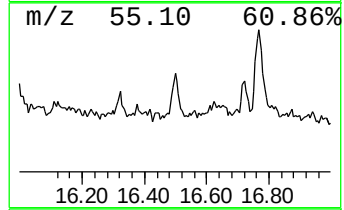
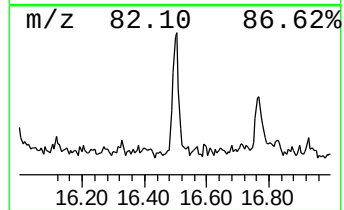
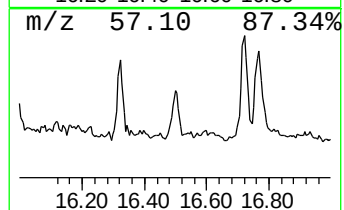
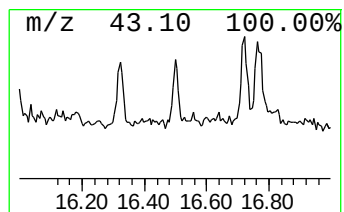
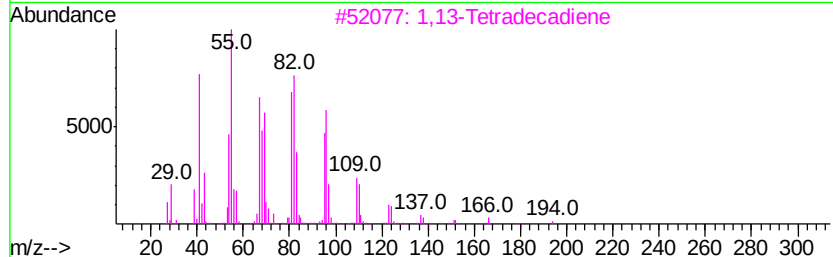
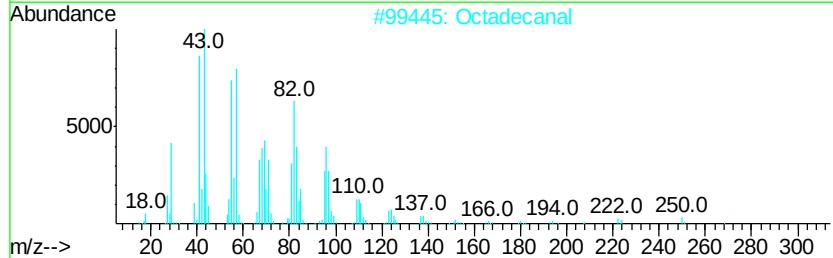
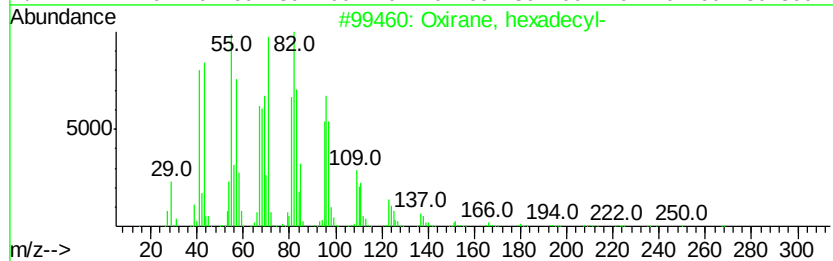
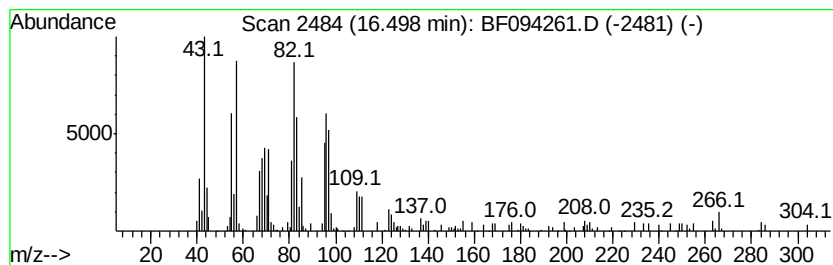
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Oxirane, hexadecyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	3.46 ng	131001	Perylene-d12	16.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Oxirane, hexadecyl-	268 C18H36O	007390-81-0 91
2		Octadecanal	268 C18H36O	000638-66-4 90
3		1,13-Tetradecadiene	194 C14H26	021964-49-8 78
4		2-Heptadecenal	252 C17H32O	1000143-48-6 78
5		11-Tetradecyn-1-ol acetate	252 C16H28O2	033925-72-3 70



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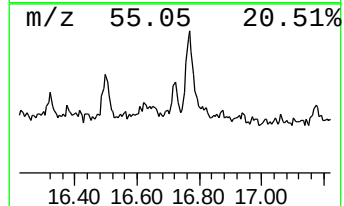
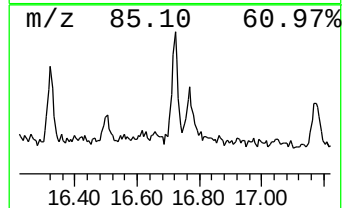
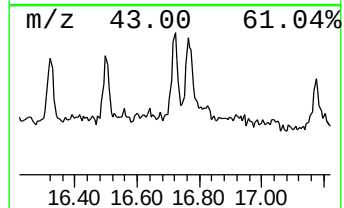
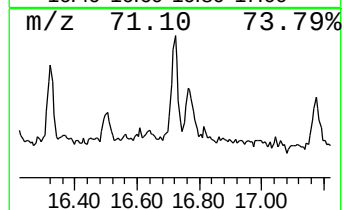
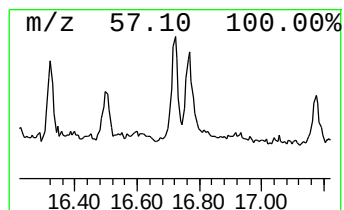
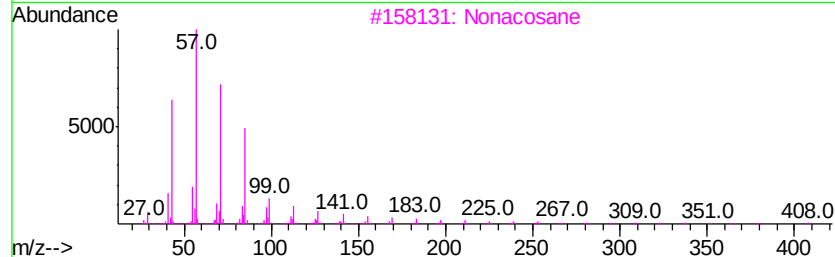
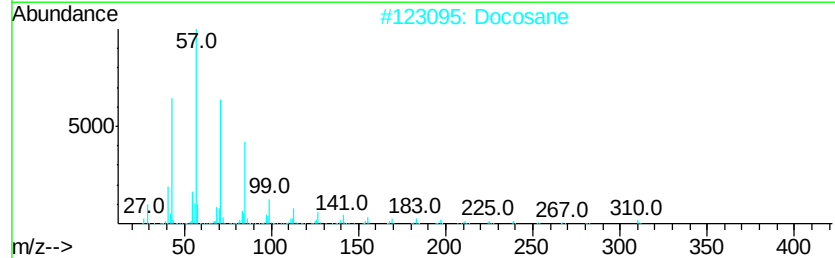
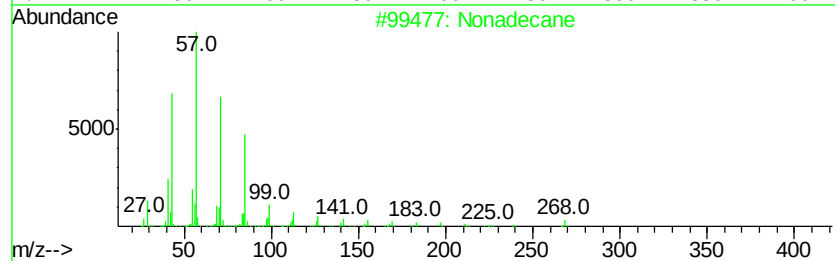
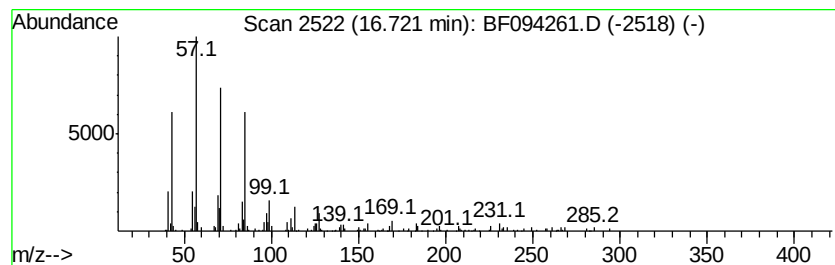
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TIC Integration Parameters: LSCINT.P

Peak Number 15 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.72	3.92 ng	148529	Perylene-d12	16.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonadecane	268	C19H40	000629-92-5	91
2		Docosane	310	C22H46	000629-97-0	90
3		Nonacosane	408	C29H60	000630-03-5	90
4		Hexadecane	226	C16H34	000544-76-3	83
5		Dodecane, 2-methyl-	184	C13H28	001560-97-0	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040617\
Data File : BF094261.D
Acq On : 7 Apr 2017 8:44
Operator : SJ/MA
Sample : I2542-18
Misc :
ALS Vial : 43 Sample Multiplier: 1

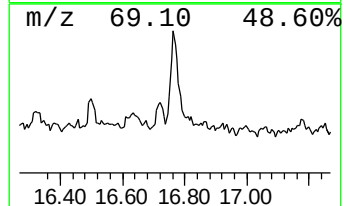
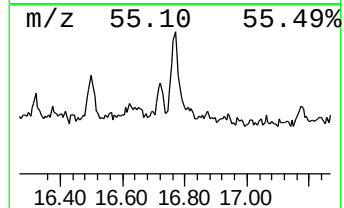
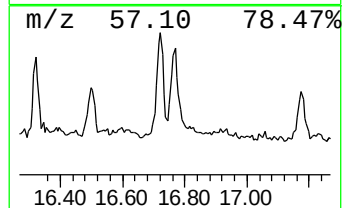
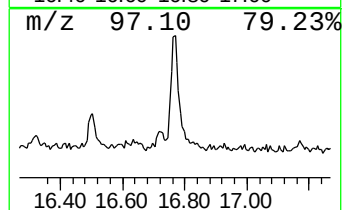
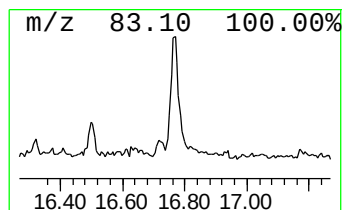
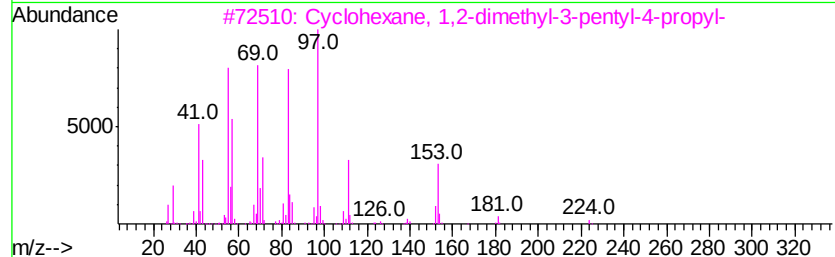
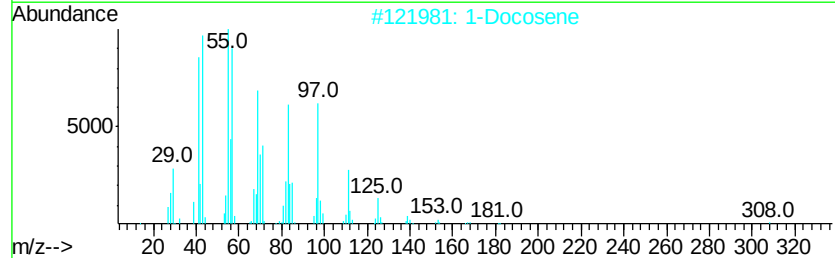
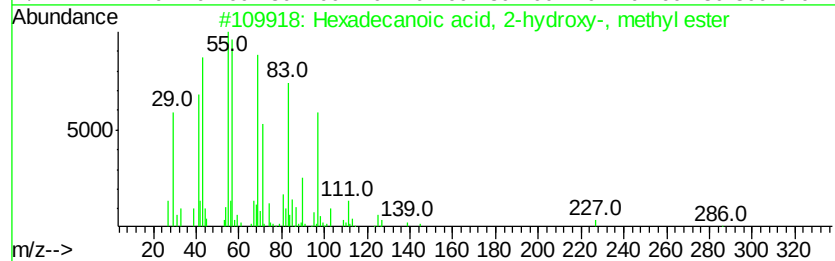
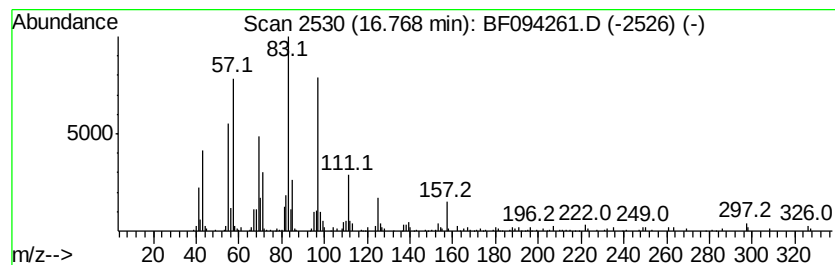
Instrument :
BNA_F
ClientSampleId :
COMP-5

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Hexadecanoic acid, 2-hydrox... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.77	6.50 ng	246102	Perylene-d12	16.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, 2-hydroxy-, m...	286	C17H34O3	016742-51-1	83
2	1-Docosene	308	C22H44	001599-67-3	55
3	Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	47
4	Cyclohexane, 1,5-diisopropyl-2,3...	196	C14H28	1000149-58-8	43
5	2-Quinuclidinone	125	C7H11NO	025289-67-2	30



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Sample : I2542-18
Misc :
ALS Vial : 43 Sample Multiplier: 1

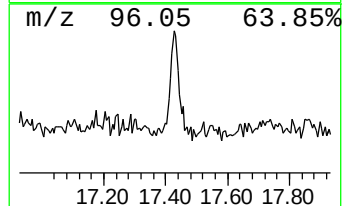
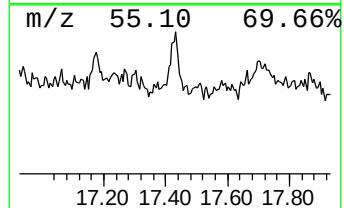
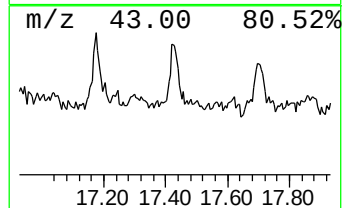
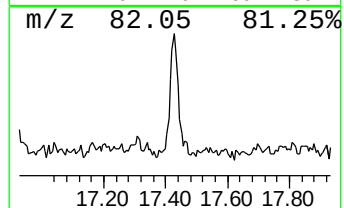
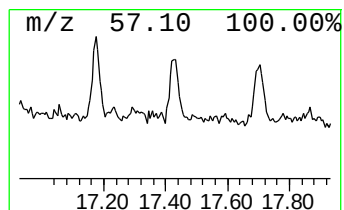
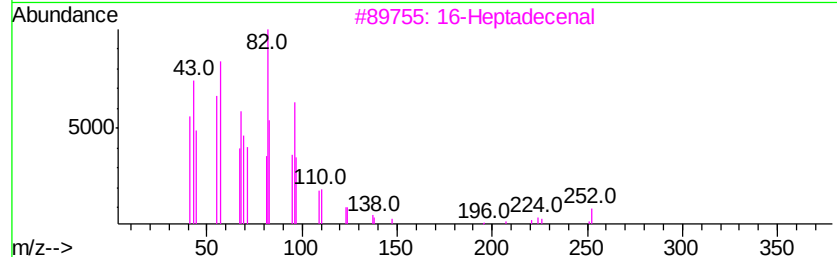
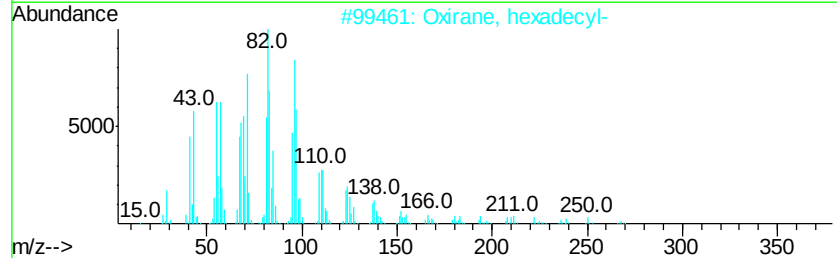
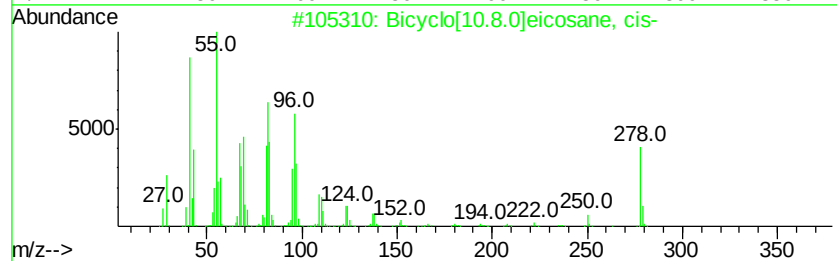
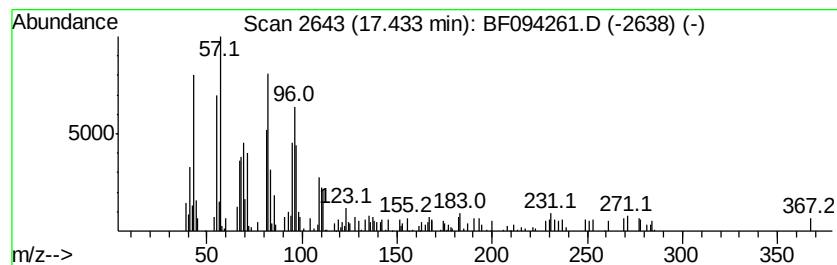
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Bicyclo[10.8.0]eicosane, cis- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.43	4.11 ng	155624	Perylene-d12	16.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	91
2	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	90
3	16-Heptadecenal	252	C17H32O	1000144-57-9	90
4	1,19-Eicosadiene	278	C20H38	014811-95-1	89
5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	87



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ALS Vial : 43 Sample Multiplier: 1

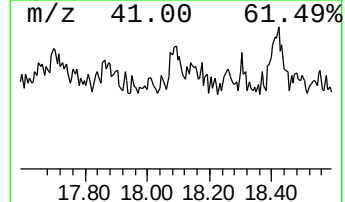
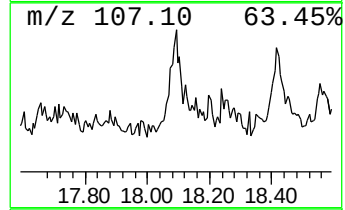
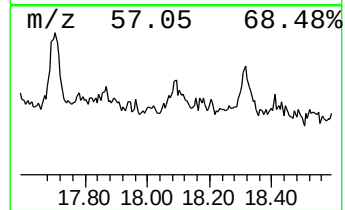
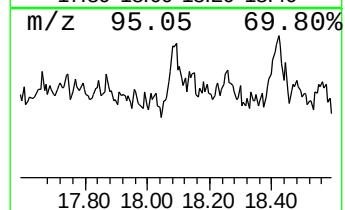
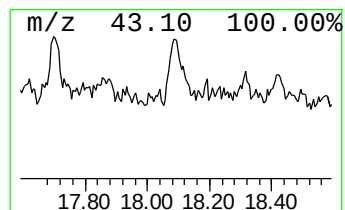
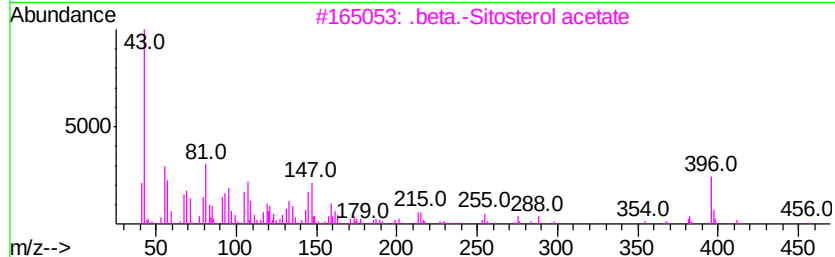
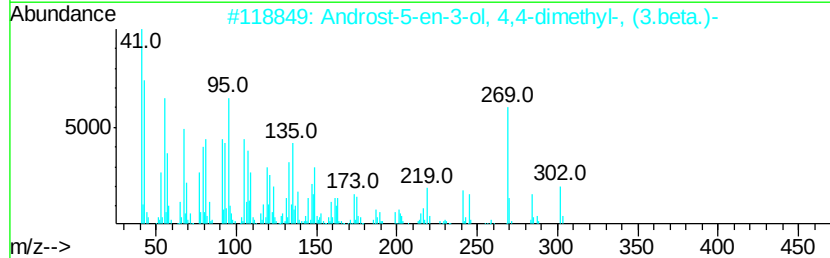
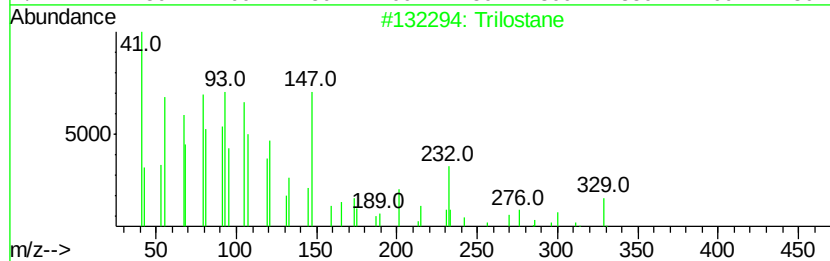
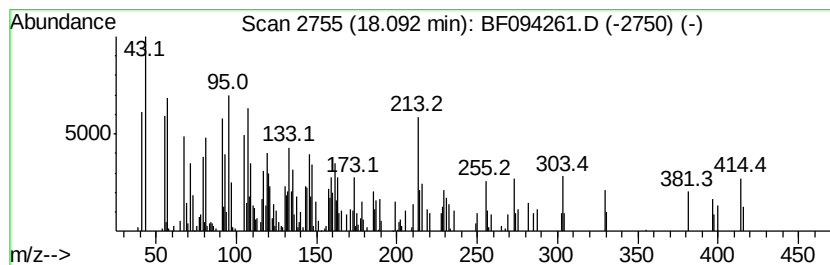
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown18.09 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.09	5.28 ng	199999	Perylene-d12	16.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trilostane	329	C20H27NO3	013647-35-3	47
2	Androst-5-en-3-ol, 4,4-dimethyl-...	302	C21H34O	007673-17-8	27
3	.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	25
4	Kauren-18-ol, acetate, (4.beta.)-	330	C22H34O2	072150-74-4	16
5	5-Cholene, 3,24-dihydroxy-	360	C24H40O2	1000251-69-2	16



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 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 COMP-5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.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...	3.99	15.7	ng	727742	1	5.85	928736	20.0
unknown5.60	5.60	74.4	ng	3455250	1	5.85	928736	20.0
Benzene, 1-methyl...	5.95	2.1	ng	95791	1	5.85	928736	20.0
Cyclopentadecane	14.47	8.9	ng	413240	5	14.57	933324	20.0
Pentadecane, 8-he...	14.87	2.1	ng	95907	5	14.57	933324	20.0
Octadecane	15.25	7.0	ng	326615	5	14.57	933324	20.0
unknown15.40	15.40	3.4	ng	127676	6	16.20	757167	20.0
Octadecane	15.61	2.5	ng	95806	6	16.20	757167	20.0
Heneicosane	15.97	12.3	ng	463773	6	16.20	757167	20.0
Pentacosane	16.32	2.1	ng	78112	6	16.20	757167	20.0
Oxirane, hexadecyl-	16.50	3.5	ng	131001	6	16.20	757167	20.0
Nonadecane	16.72	3.9	ng	148529	6	16.20	757167	20.0
Hexadecanoic acid...	16.77	6.5	ng	246102	6	16.20	757167	20.0
Bicyclo[10.8.0]ei...	17.43	4.1	ng	155624	6	16.20	757167	20.0
unknown18.09	18.09	5.3	ng	199999	6	16.20	757167	20.0