

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
 Data File : BF094702.D
 Acq On : 28 Apr 2017 23:41
 Operator : SJ/MA
 Sample : I2895-11
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-08-COMP(0-2.0)

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-BF041717.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	3.943	337	341	352	rBV	464542	586752	17.82%	1.761%
2	4.331	403	407	421	rBV	3164574	3189076	96.83%	9.570%
3	4.713	469	472	478	rBV	37621	40166	1.22%	0.121%
4	5.401	584	589	592	rBV	2754787	3169211	96.23%	9.511%
5	5.525	605	610	613	rBV	3269270	3293372	100.00%	9.883%
6	5.772	648	652	656	rBV	697721	660321	20.05%	1.982%
7	5.948	677	682	685	rBV	2620116	2426768	73.69%	7.283%
8	6.419	757	762	765	rBV	1785647	1987135	60.34%	5.963%
9	7.260	901	905	911	rVB	988861	865308	26.27%	2.597%
10	7.819	994	1000	1005	rBV2	49988	53506	1.62%	0.161%
11	8.554	1122	1125	1126	rBV	41798	38276	1.16%	0.115%
12	8.583	1126	1130	1133	rVB	3022635	2818999	85.60%	8.460%
13	9.083	1212	1215	1219	rBV	249146	232729	7.07%	0.698%
14	9.119	1219	1221	1224	rVB	42893	36233	1.10%	0.109%
15	9.395	1264	1268	1273	rVB	868012	820134	24.90%	2.461%
16	9.730	1316	1325	1328	rBV6	15043	33750	1.02%	0.101%
17	10.372	1429	1434	1442	rBV	1442255	1585060	48.13%	4.757%
18	10.577	1466	1469	1470	rBV	34924	35704	1.08%	0.107%
19	11.219	1574	1578	1581	rBV	736952	717377	21.78%	2.153%
20	11.242	1581	1582	1587	rVB	85950	76111	2.31%	0.228%
21	11.954	1699	1703	1705	rBV	91759	101238	3.07%	0.304%
22	12.713	1828	1832	1837	rBV	213383	234954	7.13%	0.705%
23	12.918	1865	1867	1874	rBV8	22198	36076	1.10%	0.108%
24	12.983	1874	1878	1882	rVV	222794	234517	7.12%	0.704%
25	13.030	1882	1886	1891	rVB5	40862	56506	1.72%	0.170%
26	13.195	1909	1914	1918	rBV	1769434	1916897	58.20%	5.752%
27	13.401	1947	1949	1954	rVB2	41778	49764	1.51%	0.149%
28	13.542	1968	1973	1977	rVB5	52816	77470	2.35%	0.232%
29	14.148	2073	2076	2080	rBV2	94442	107815	3.27%	0.324%
30	14.265	2092	2096	2100	rBV2	208810	225798	6.86%	0.678%
31	14.371	2109	2114	2119	rVB3	250992	319807	9.71%	0.960%
32	14.424	2119	2123	2126	rBV2	67329	78263	2.38%	0.235%
33	14.471	2126	2131	2134	rVV2	608374	756673	22.98%	2.271%
34	14.524	2134	2140	2145	rVB	1964442	2148286	65.23%	6.447%

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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	14.871	2196	2199	2202	rBV2	36916	42778	1.30%	0.128%
36	15.036	2224	2227	2231	rBV2	63286	65845	2.00%	0.198%
37	15.142	2241	2245	2254	rBV	966523	1118886	33.97%	3.358%
38	15.277	2265	2268	2274	rVB	66734	110285	3.35%	0.331%
39	15.659	2330	2333	2336	rBV2	84007	85786	2.60%	0.257%
40	15.701	2336	2340	2343	rBV2	106911	155495	4.72%	0.467%
41	15.859	2363	2367	2376	rBV	744174	946895	28.75%	2.842%
42	15.983	2385	2388	2392	rBV2	108070	133799	4.06%	0.402%
43	16.042	2396	2398	2402	rBV	67924	77071	2.34%	0.231%
44	16.101	2404	2408	2411	rBV	429805	450447	13.68%	1.352%
45	16.377	2451	2455	2467	rVB3	190295	324182	9.84%	0.973%
46	16.589	2487	2491	2496	rVB	139950	182218	5.53%	0.547%
47	17.265	2601	2606	2613	rVB	196836	335904	10.20%	1.008%
48	18.342	2783	2789	2799	rVB10	104135	283560	8.61%	0.851%

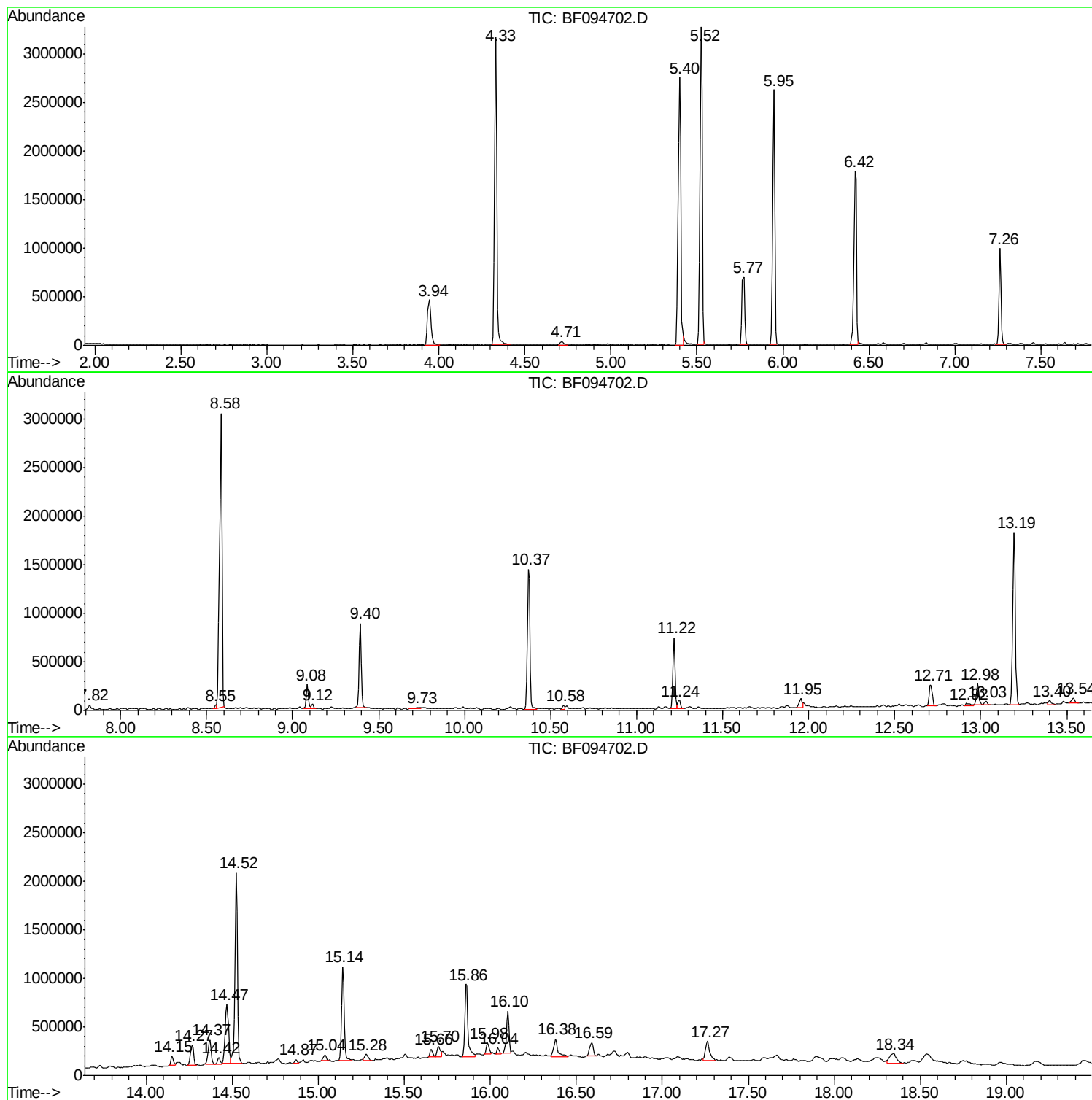
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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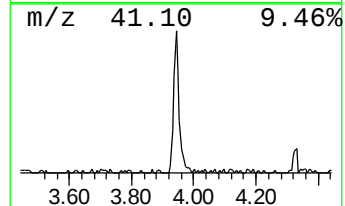
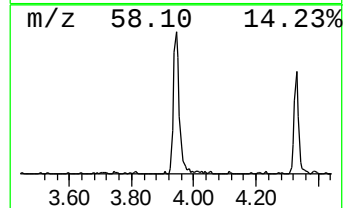
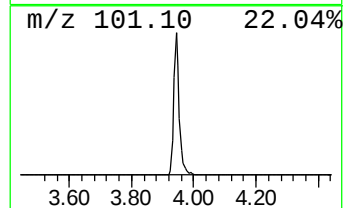
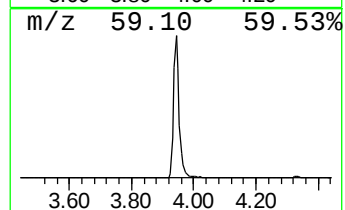
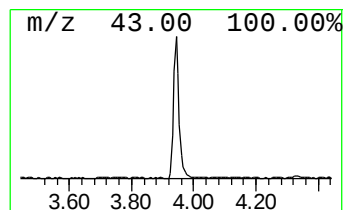
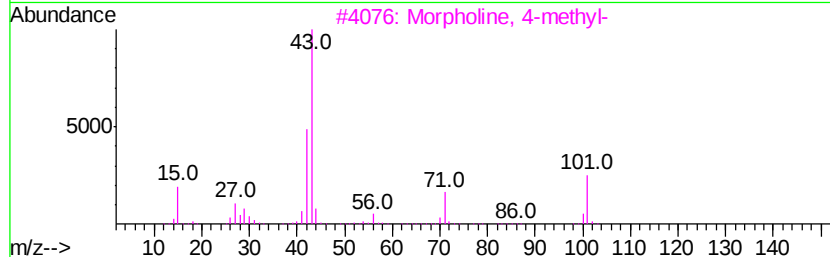
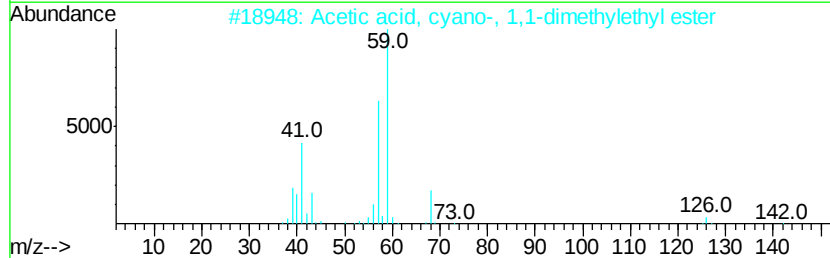
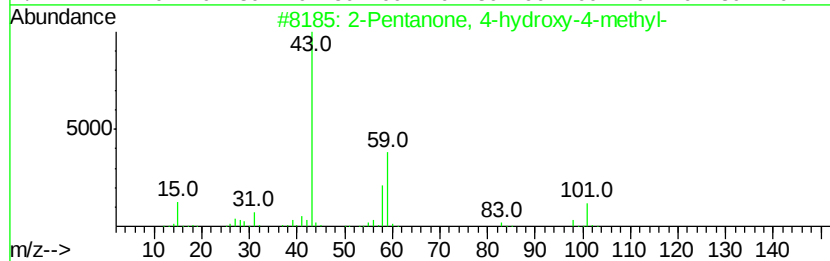
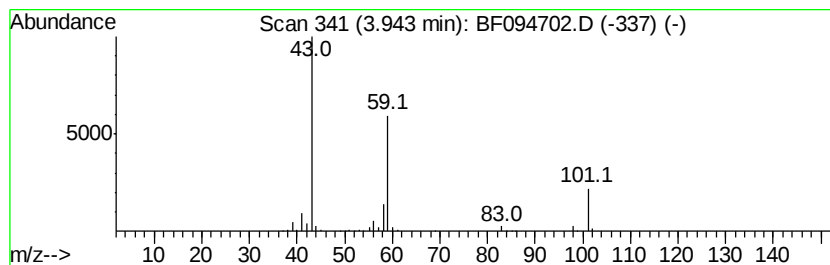
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.94	17.77 ng	586752	1,4-Dichlorobenzene-d4	5.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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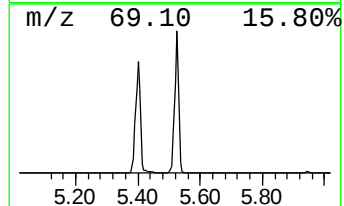
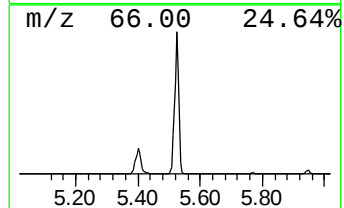
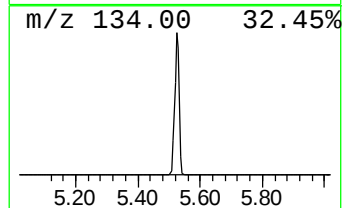
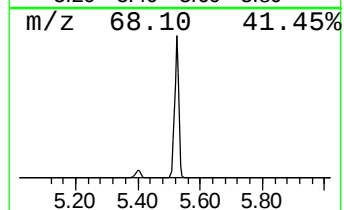
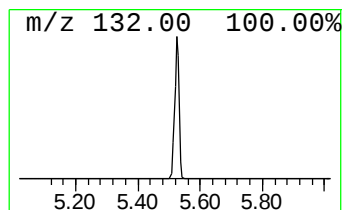
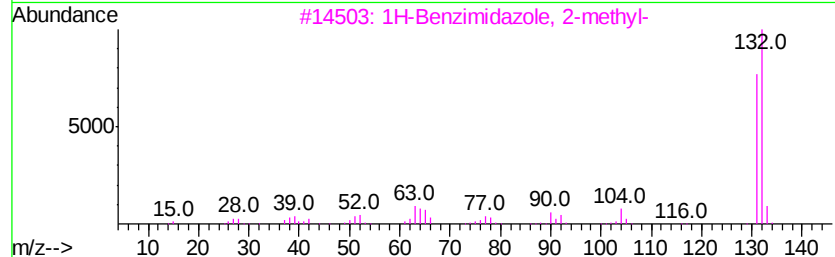
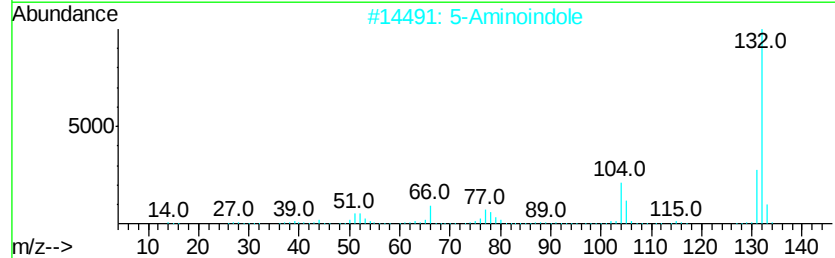
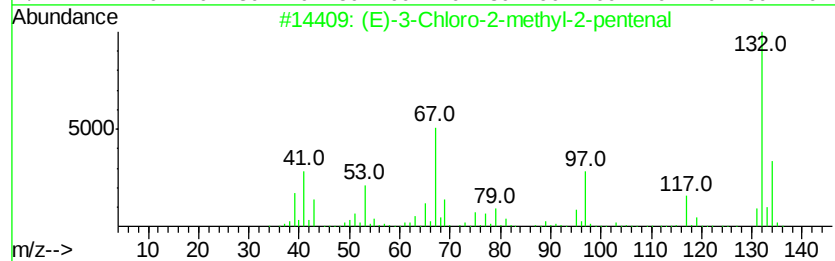
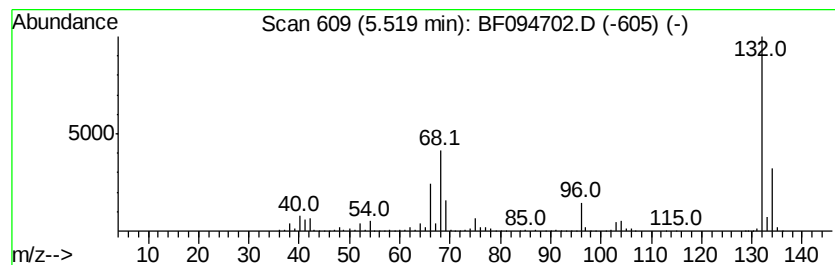
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown5.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.52	99.75 ng	3293370	1,4-Dichlorobenzene-d4	5.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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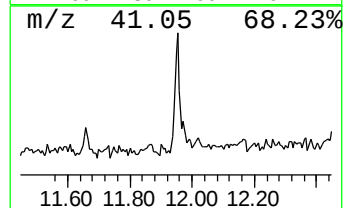
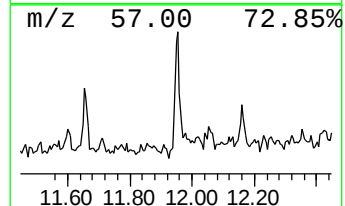
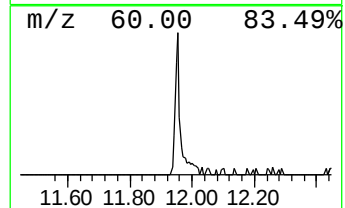
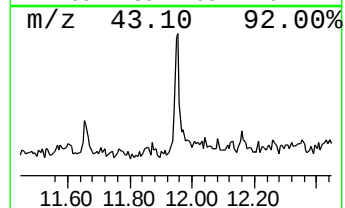
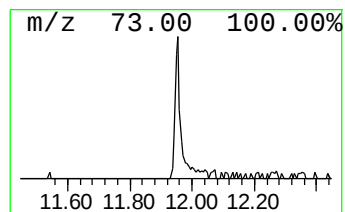
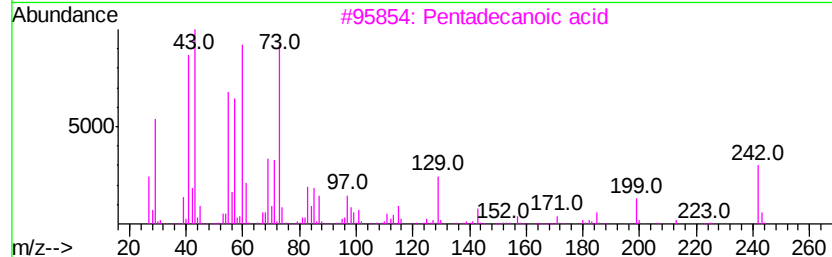
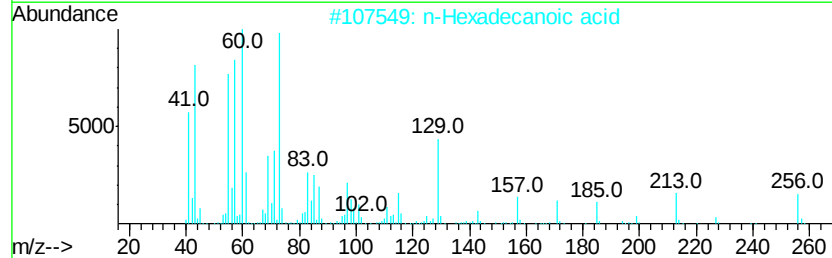
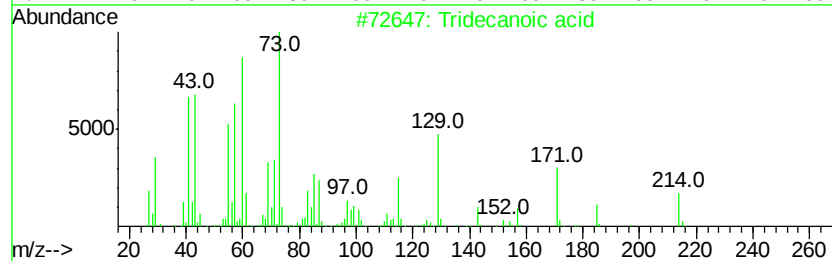
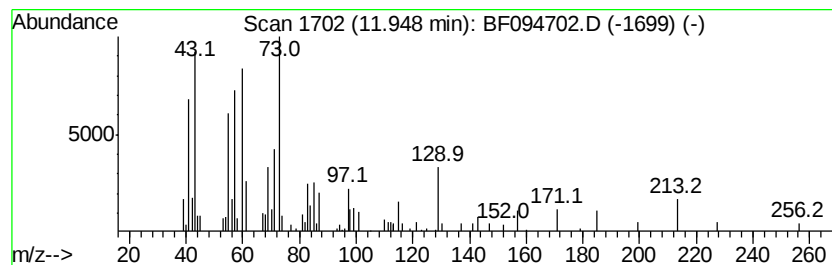
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Tridecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.95	2.82 ng	101238	Phenanthrene-d10	11.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecanoic acid	214	C13H26O2	000638-53-9	95
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	80
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		Dodecanoic acid	200	C12H24O2	000143-07-7	64



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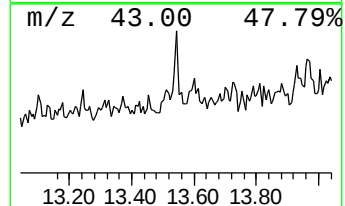
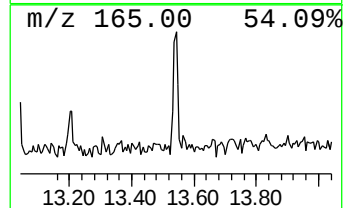
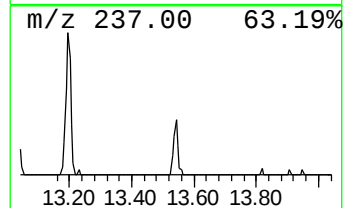
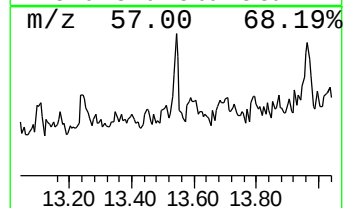
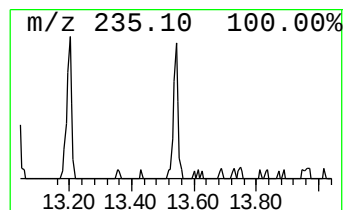
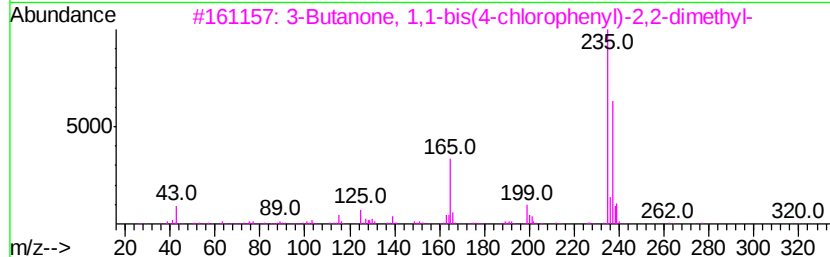
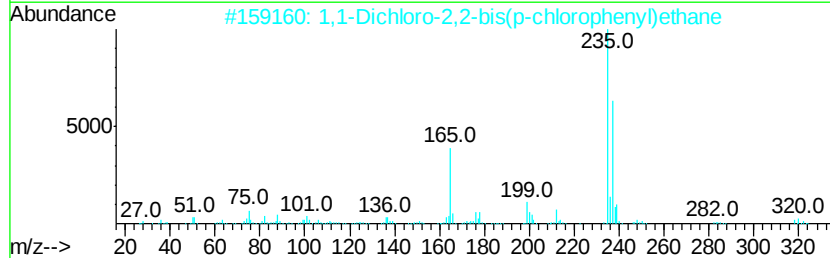
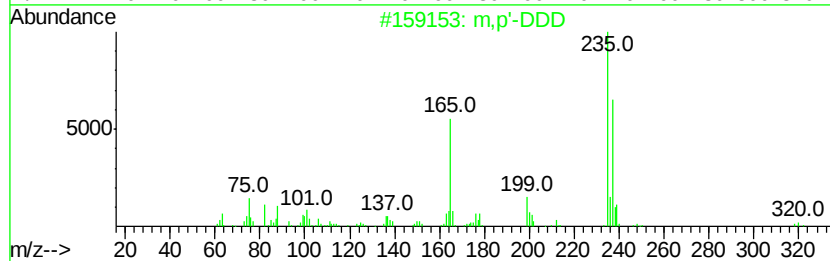
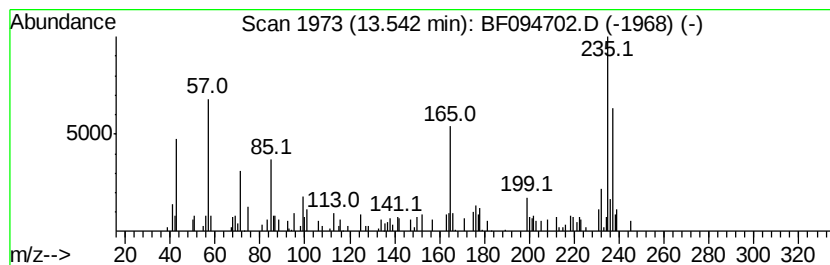
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 m,p'-DDD Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	2.05 ng	77470	Chrysene-d12	14.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	m,p'-DDD	318	C14H10Cl4	004329-12-8	64
2		1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	64
3		3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	58
4		1,2-Bis(2-chlorophenyl)-1,2-bis(...	470	C26H18Cl4	1000137-94-8	53
5		9H-Carbazole, 3,6-dichloro-	235	C12H7Cl2N	005599-71-3	49



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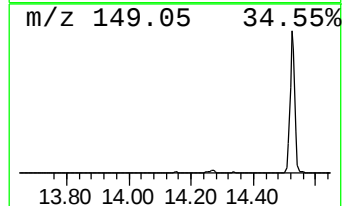
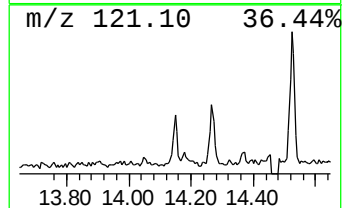
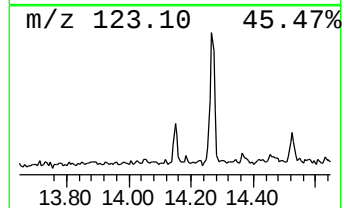
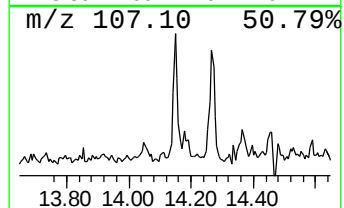
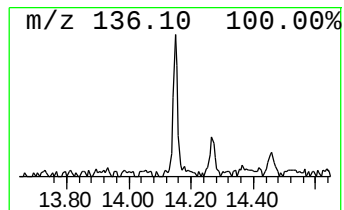
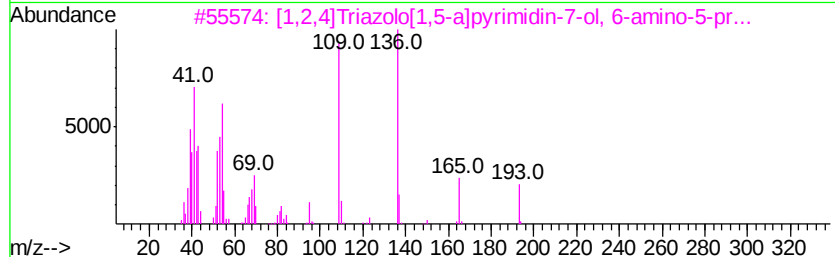
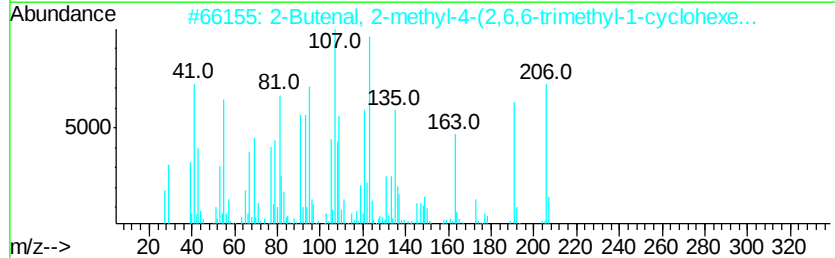
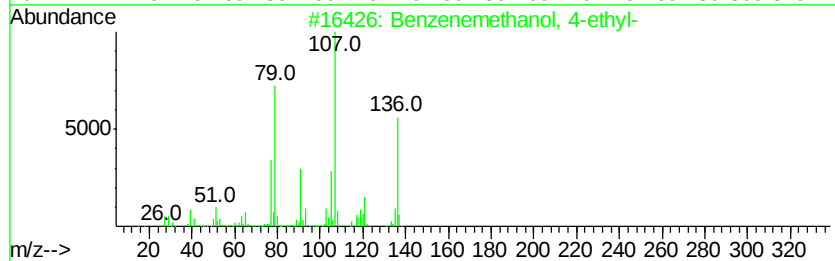
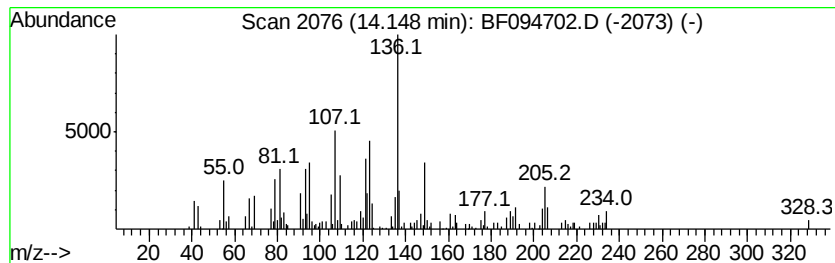
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown14.15 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.15	2.85 ng	107815	Chrysene-d12	14.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenemethanol, 4-ethyl-	136	C9H12O	000768-59-2	46
2			2-Butenal, 2-methyl-4-(2,6,6-tri...	206	C14H22O	003155-71-3	38
3			[1,2,4]Triazolo[1,5-a]pyrimidin-...	193	C8H11N5O	1000362-66-7	35
4			Glutaric acid, monoamide, N-(2-m...	307	C17H25NO4	1000360-02-5	27
5			Benzenemethanol, 3,5-dimethyl-	136	C9H12O	027129-87-9	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042817\
Data File : BF094702.D
Acq On : 28 Apr 2017 23:41
Operator : SJ/MA
Sample : I2895-11
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-08-COMP(0-2.0)

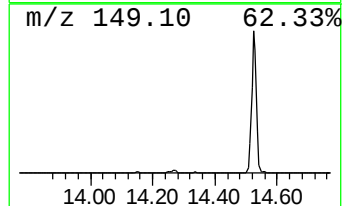
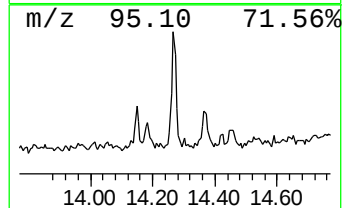
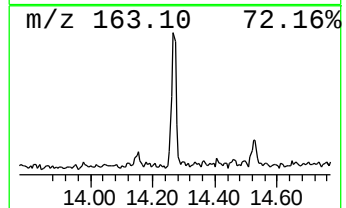
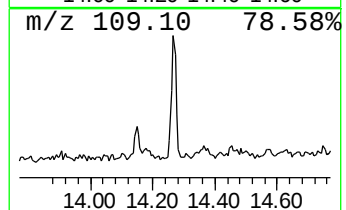
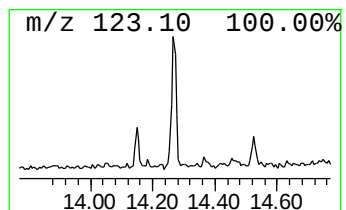
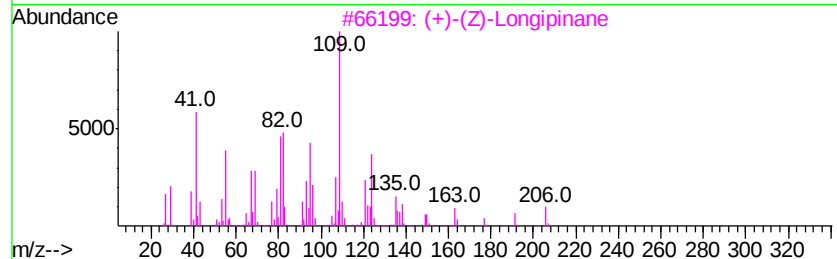
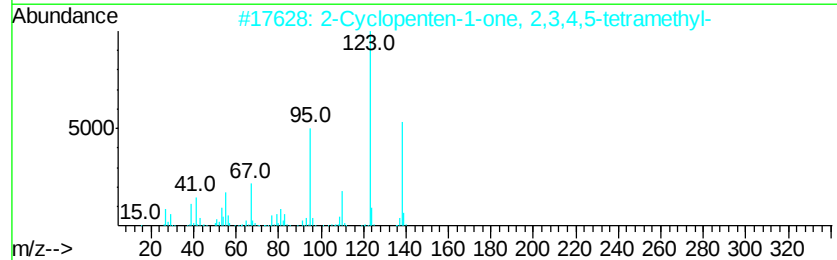
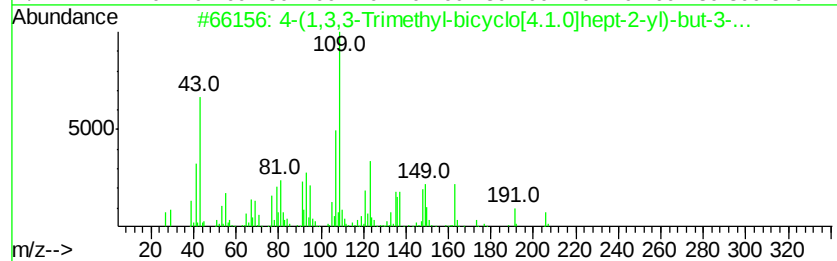
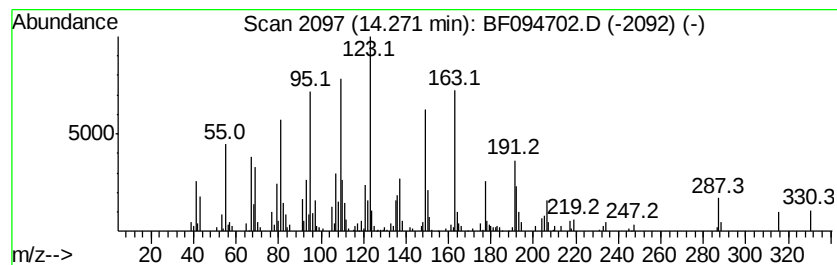
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown14.27 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	5.97 ng	225798	Chrysene-d12	14.47

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-	(1,3,3-Trimethyl-bicyclo[4.1.0...	206	C14H22O	077143-31-8	41	
2	2-	Cyclopenten-1-one, 2,3,4,5-tet...	138	C9H14O	054458-61-6	38	
3	(+)-(Z)-	Longipinane	206	C15H26	1000156-12-3	35	
4	1-(1-Ethyl-2,3-dimethyl-cyclopen...	166	C11H18O	1000185-18-6	30		
5	Cyclohexene, 3-methyl-6-(1-methy...	138	C10H18	005256-65-5	25		



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
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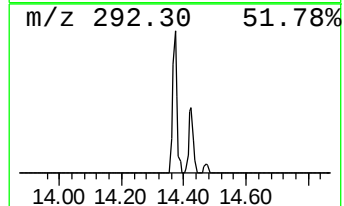
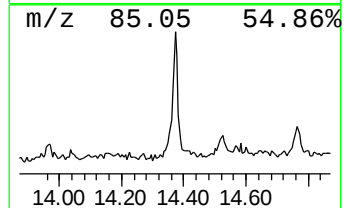
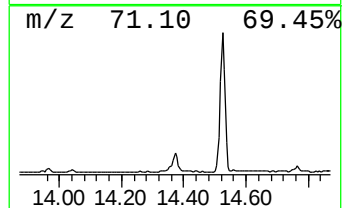
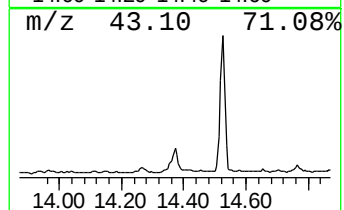
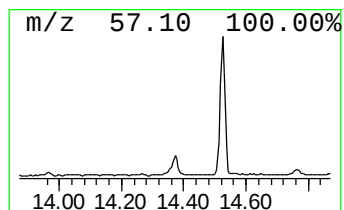
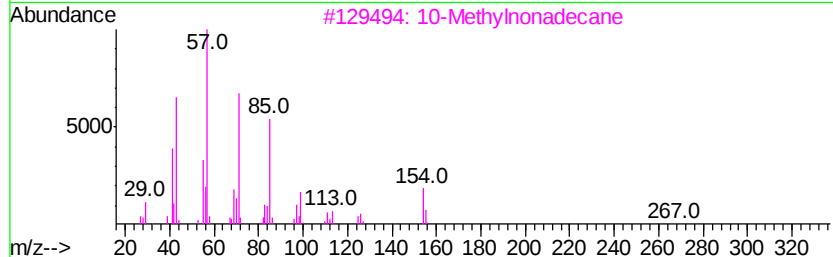
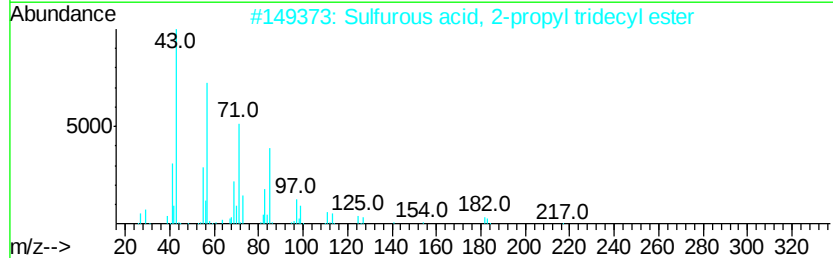
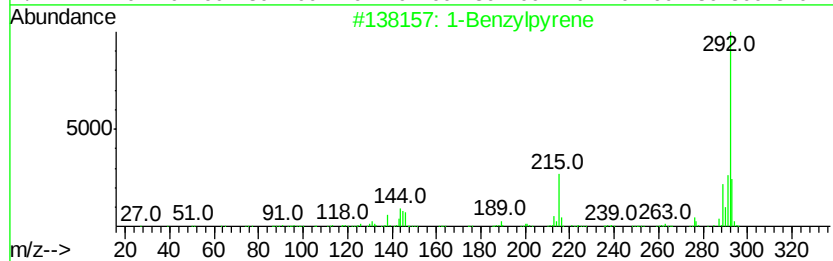
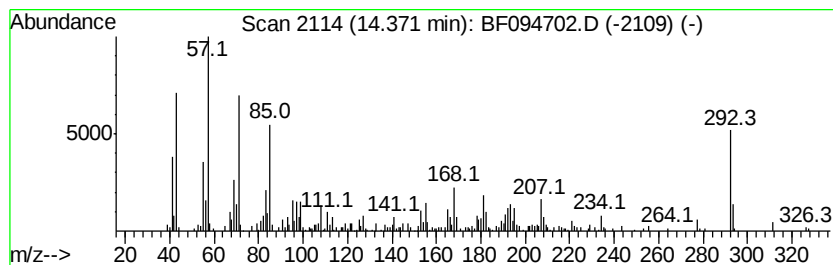
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown14.37 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.37	8.45 ng	319807	Chrysene-d12	14.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Benzylpyrene	292	C23H16	042211-34-7	43
2		Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	41
3		10-Methylnonadecane	282	C20H42	056862-62-5	41
4		Hentriacontane	437	C31H64	000630-04-6	41
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	41



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
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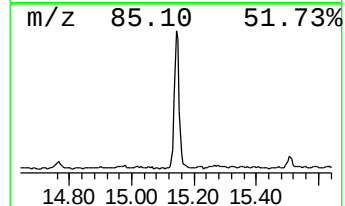
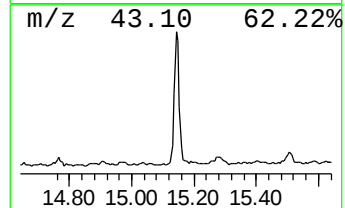
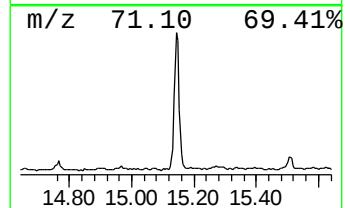
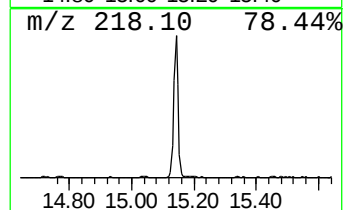
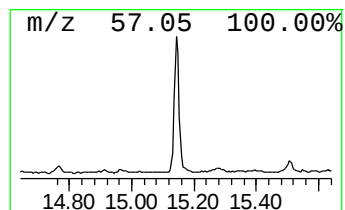
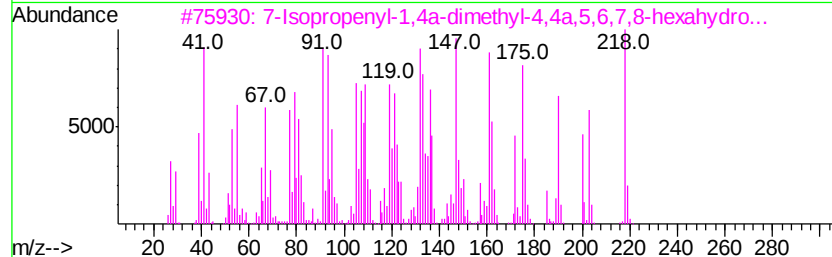
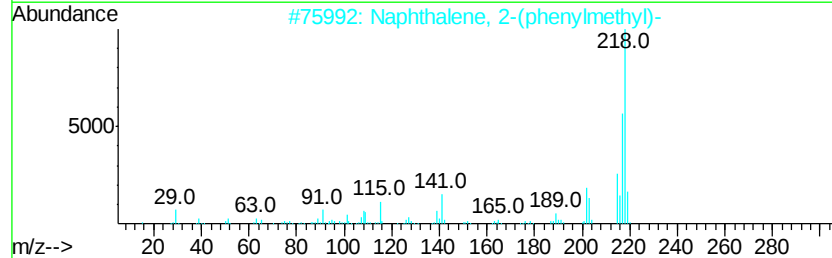
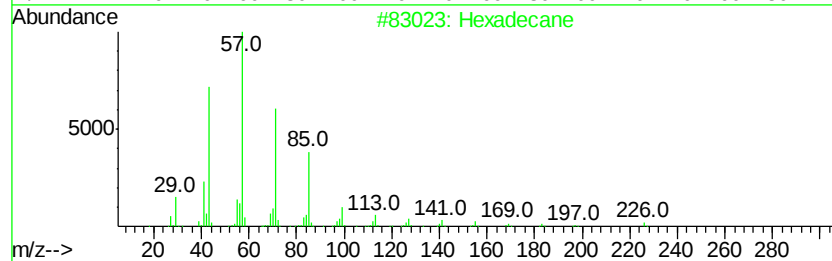
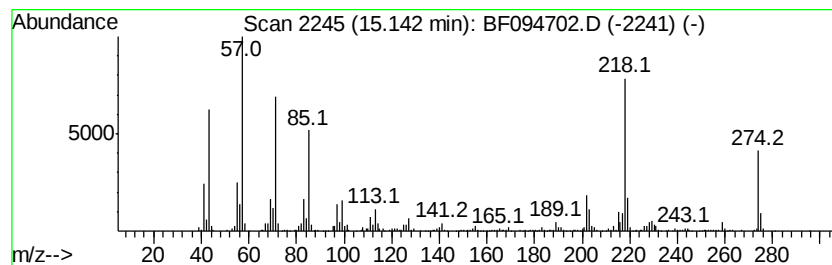
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	29.57 ng	1118890	Chrysene-d12	14.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Naphthalene, 2-(phenylmethyl)-	218	C17H14	000613-59-2	38
3		7-Isopropenyl-1,4a-dimethyl-4,4a...	218	C15H22O	000473-08-5	25
4		Heptadecane	240	C17H36	000629-78-7	20
5		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	20



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
Data File : BF094702.D
Acq On : 28 Apr 2017 23:41
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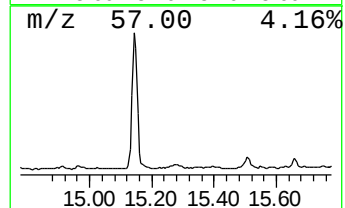
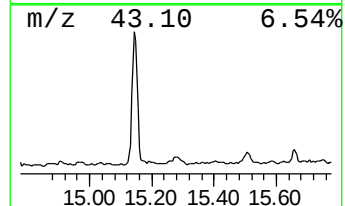
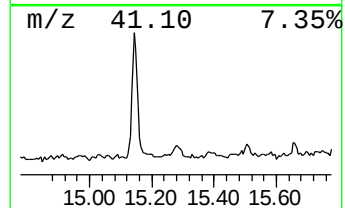
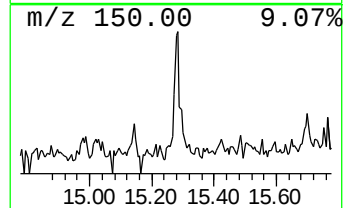
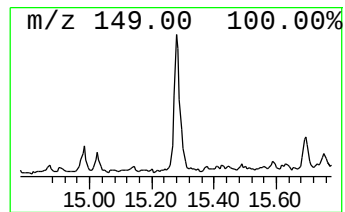
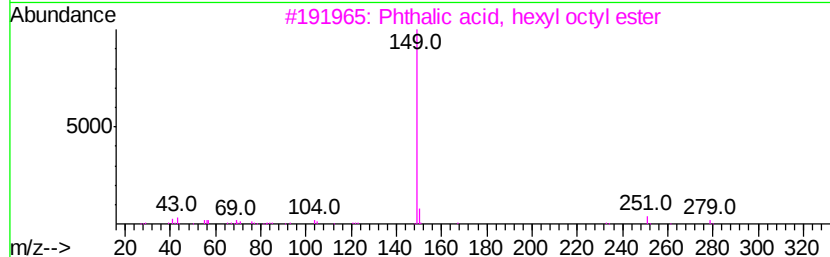
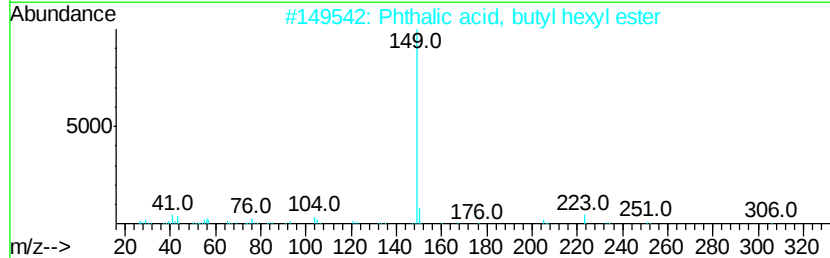
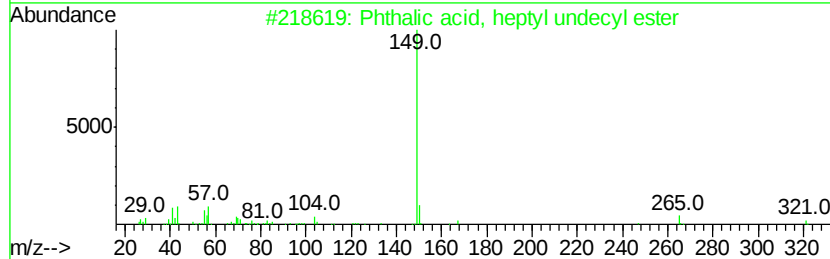
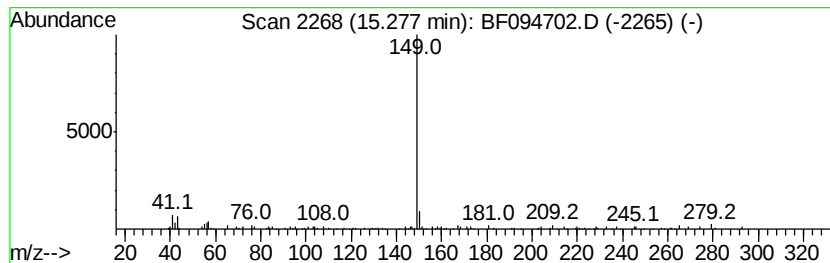
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phthalic acid, heptyl undec... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.28	2.92 ng	110285	Chrysene-d12	14.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic acid, heptyl undecyl ester	418	C26H42O4	1000308-94-0	72
2		Phthalic acid, butyl hexyl ester	306	C18H26O4	1000308-99-5	72
3		Phthalic acid, hexyl octyl ester	362	C22H34O4	1000308-97-8	72
4		Phthalic acid, di(hex-3-yl) ester	334	C20H30O4	1000356-96-7	72
5		Phthalic acid, 2-chloropropyl no...	368	C20H29ClO4	1000356-83-0	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
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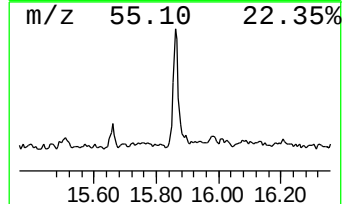
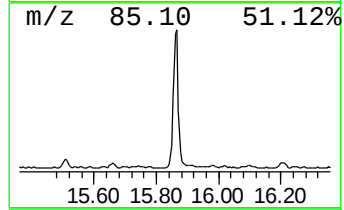
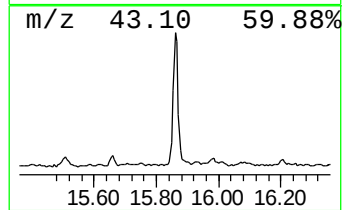
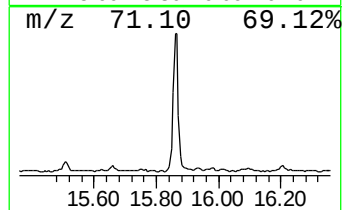
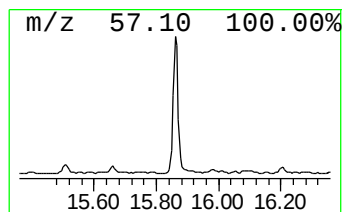
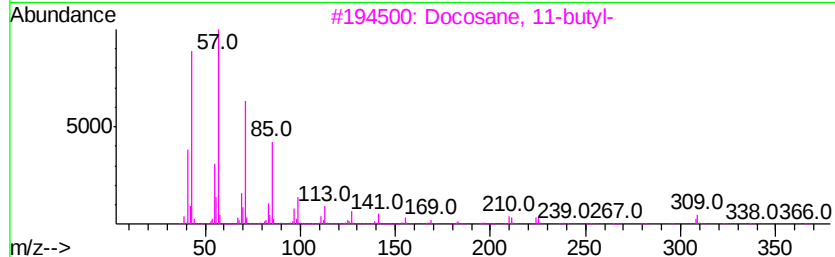
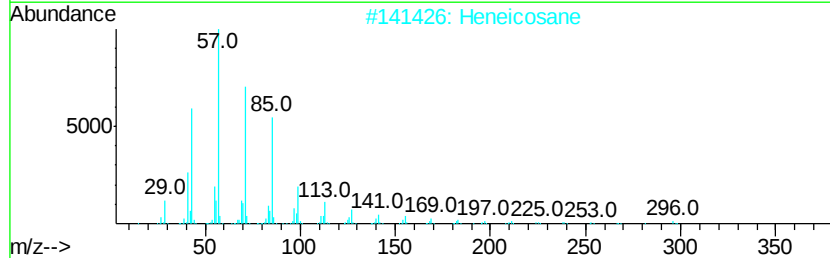
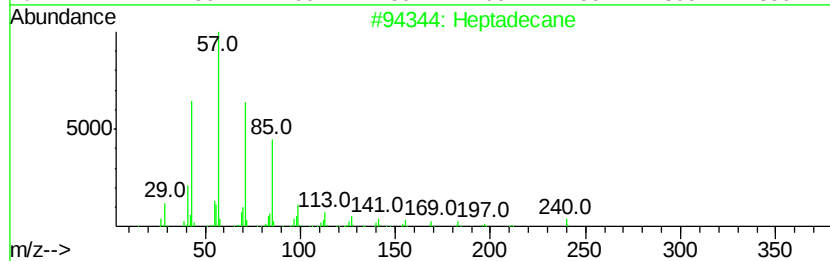
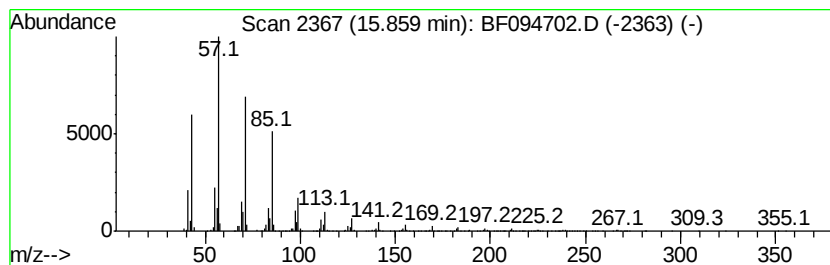
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	42.04 ng	946895	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	93
2		Heneicosane	296	C21H44	000629-94-7	91
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



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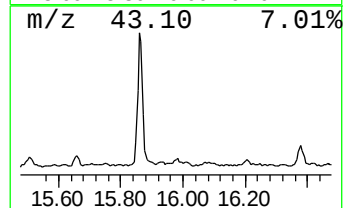
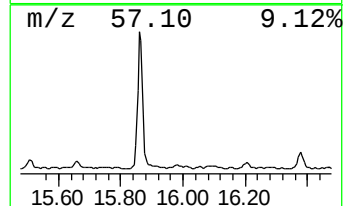
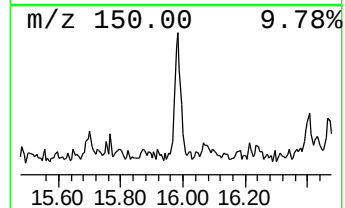
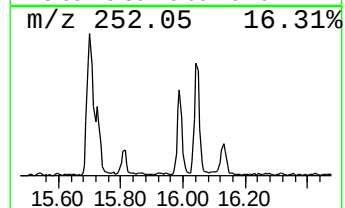
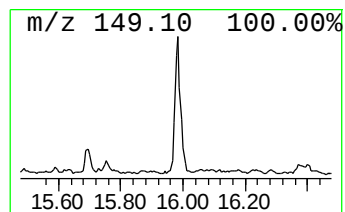
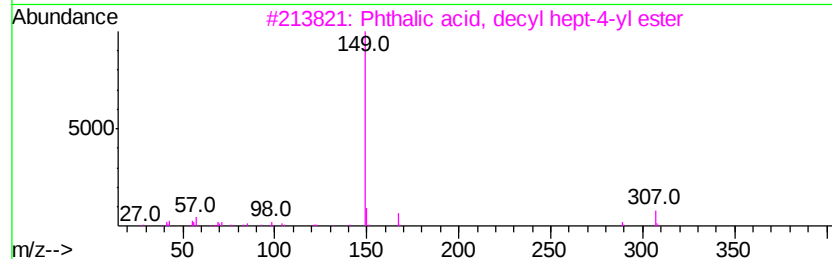
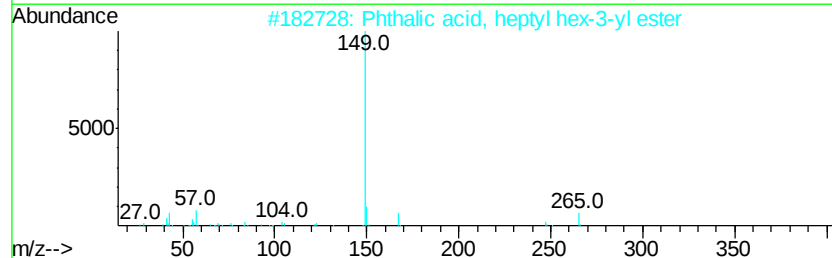
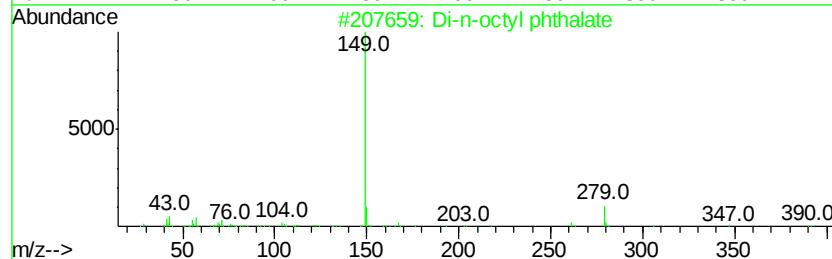
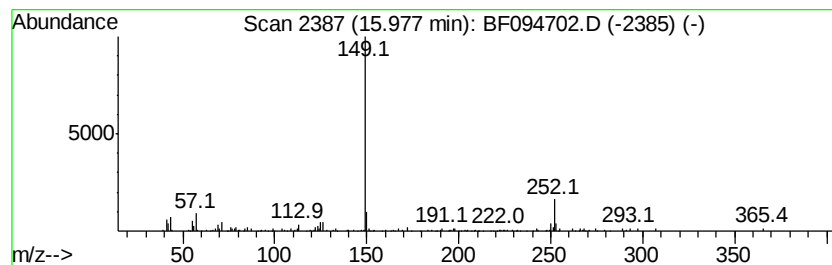
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown15.98 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.98	5.94 ng	133799	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Di-n-octyl phthalate	390	C24H38O4	000117-84-0	46
2		Phthalic acid, heptyl hex-3-yl e...	348	C21H32O4	1000356-95-8	43
3		Phthalic acid, decyl hept-4-yl e...	404	C25H40O4	1000356-79-1	43
4		Phthalic acid, 2-chloropropyl he...	480	C28H45ClO4	1000356-83-8	43
5		Phthalic acid, nonyl tridec-2-yn...	470	C30H46O4	1000315-44-2	43



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ALS Vial : 23 Sample Multiplier: 1

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ClientSampleId :
SB-08-COMP(0-2.0)

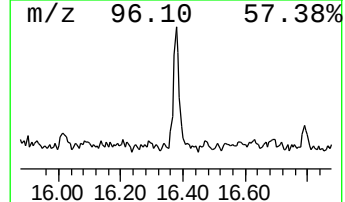
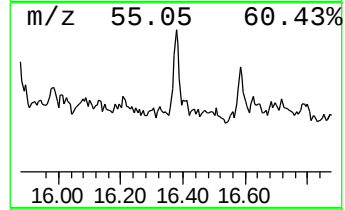
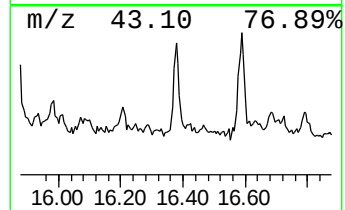
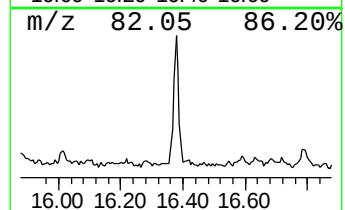
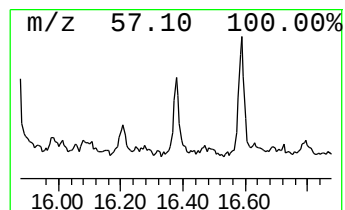
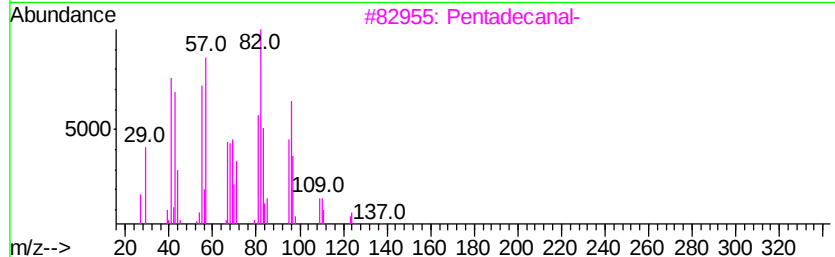
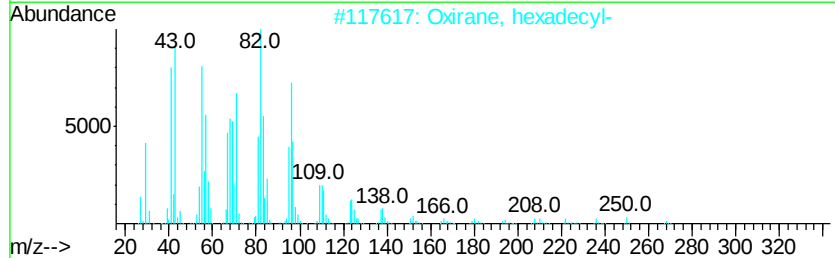
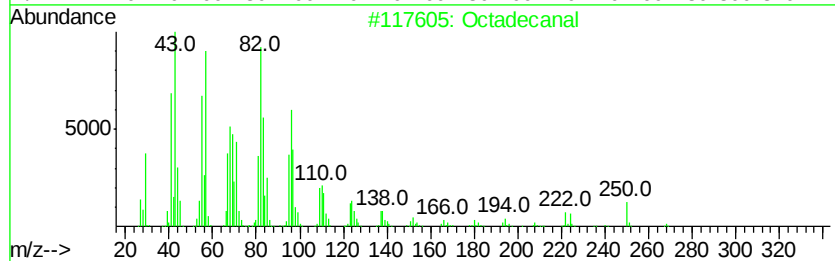
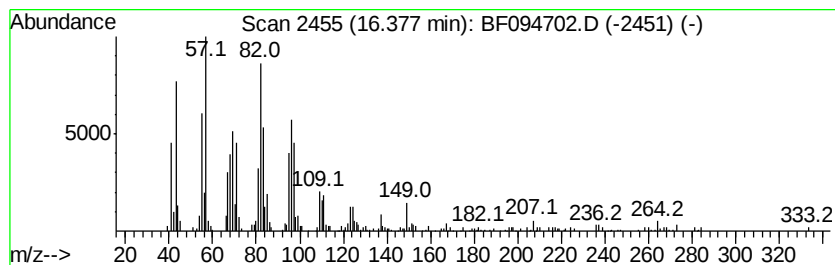
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecanal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.38	14.39 ng	324182	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanal	268	C18H36O	000638-66-4	95
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
3		Pentadecanal-	226	C15H30O	002765-11-9	90
4		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	81
5		Hexadecanal	240	C16H32O	000629-80-1	68



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042817\
Data File : BF094702.D
Acq On : 28 Apr 2017 23:41
Operator : SJ/MA
Sample : I2895-11
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-08-COMP(0-2.0)

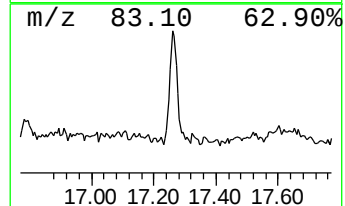
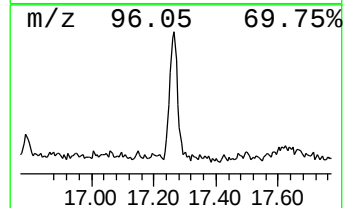
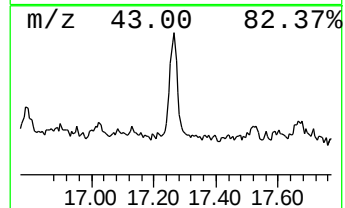
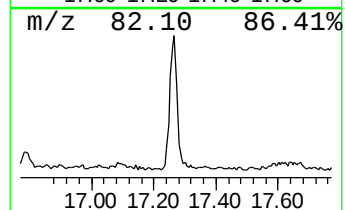
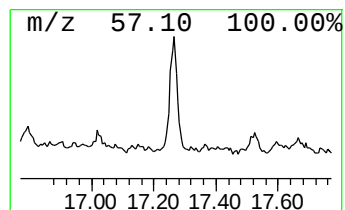
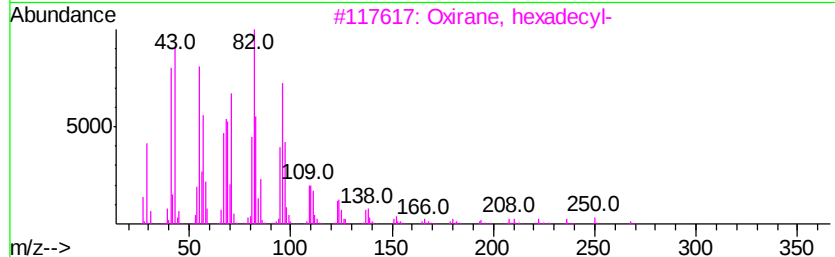
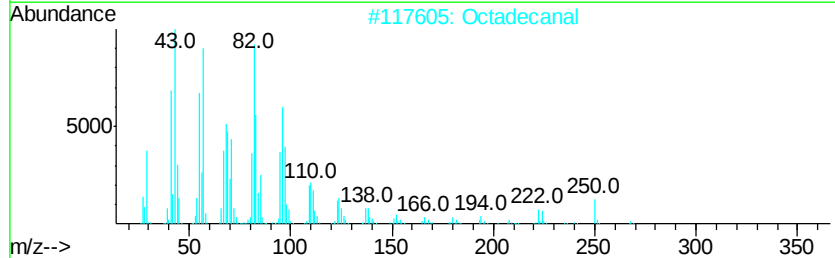
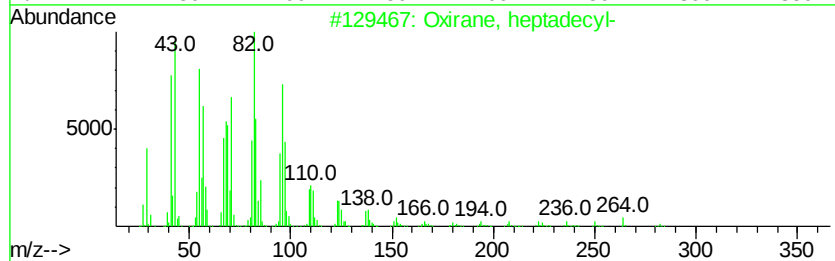
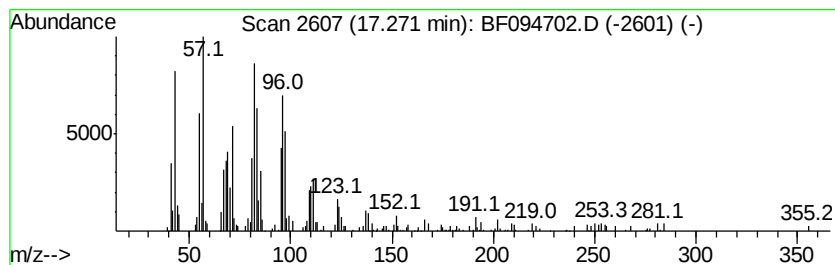
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Oxirane, heptadecyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.27	14.91 ng	335904	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	94
2		Octadecanal	268	C18H36O	000638-66-4	91
3		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	87
4		Pentadecanal-	226	C15H30O	002765-11-9	86
5		Tetradecanal	212	C14H28O	000124-25-4	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042817\
Data File : BF094702.D
Acq On : 28 Apr 2017 23:41
Operator : SJ/MA
Sample : I2895-11
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-08-COMP(0-2.0)

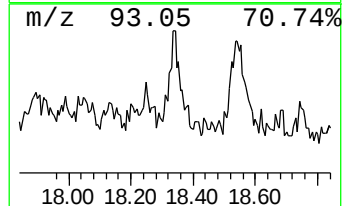
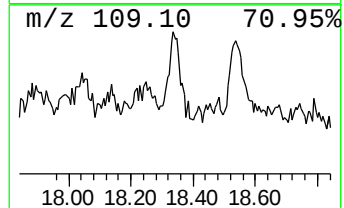
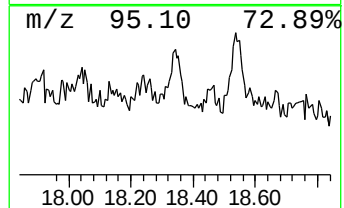
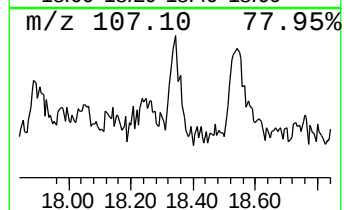
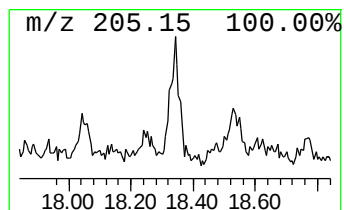
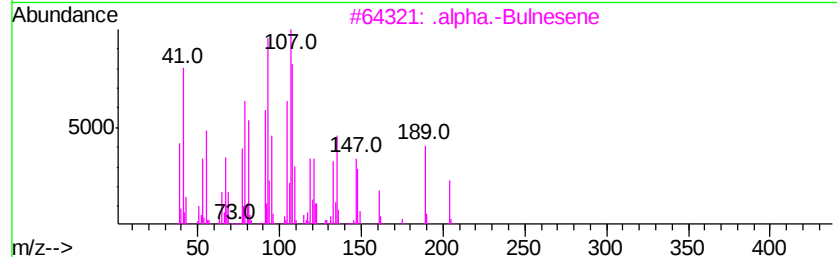
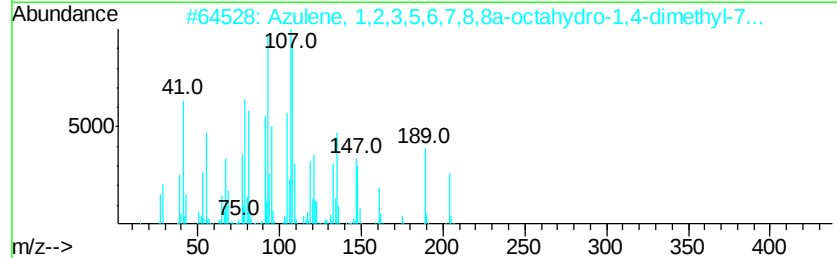
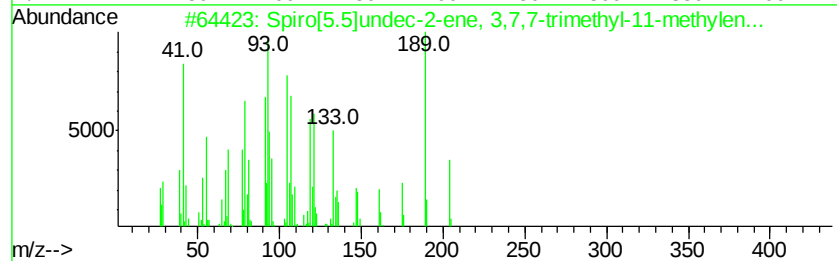
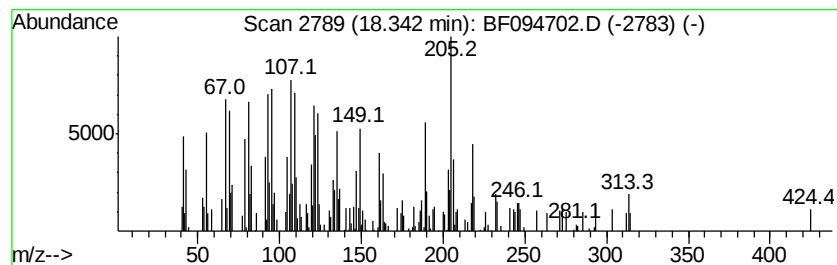
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Spiro[5.5]undec-2-ene, 3,7,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.34	12.59 ng	283560	Perylene-d12	16.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Spiro[5.5]undec-2-ene, 3,7,7-tri...	204	C15H24	018431-82-8	53
2		Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	47
3		.alpha.-Bulnesene	204	C15H24	1000374-19-9	47
4		Guaia-9,11-diene	204	C15H24	1000374-19-8	45
5		2-Hydroxymethyl-2,6,8,8-tetramet...	236	C16H28O	137235-47-3	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042817\
 Data File : BF094702.D
 Acq On : 28 Apr 2017 23:41
 Operator : SJ/MA
 Sample : I2895-11
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-08-COMP(0-2.0)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	3.94	17.8	ng	586752	1	5.77	660321	20.0
unknown5.52	5.52	99.8	ng	3293370	1	5.77	660321	20.0
Tridecanoic acid	11.95	2.8	ng	101238	4	11.22	717377	20.0
m,p'-DDD	13.54	2.0	ng	77470	5	14.47	756673	20.0
unknown14.15	14.15	2.9	ng	107815	5	14.47	756673	20.0
unknown14.27	14.27	6.0	ng	225798	5	14.47	756673	20.0
unknown14.37	14.37	8.4	ng	319807	5	14.47	756673	20.0
Hexadecane	15.14	29.6	ng	1118890	5	14.47	756673	20.0
Phthalic acid, he...	15.28	2.9	ng	110285	5	14.47	756673	20.0
Heptadecane	15.86	42.0	ng	946895	6	16.10	450447	20.0
unknown15.98	15.98	5.9	ng	133799	6	16.10	450447	20.0
Octadecanal	16.38	14.4	ng	324182	6	16.10	450447	20.0
Oxirane, heptadecyl-	17.27	14.9	ng	335904	6	16.10	450447	20.0
Spiro[5.5]undec-2...	18.34	12.6	ng	283560	6	16.10	450447	20.0