

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022118\
 Data File : BF103132.D
 Acq On : 21 Feb 2018 13:24
 Operator : SJ/JU
 Sample : J1404-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-022018

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-BF020318.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	2.134	3	8	19	rVB	854748	1099163	9.35%	1.827%
2	5.498	576	580	590	rBV	3295357	3411850	29.01%	5.670%
3	6.510	747	752	755	rBV	3111988	3118012	26.52%	5.181%
4	6.651	771	776	779	rBV	3246564	3287819	27.96%	5.464%
5	6.875	811	814	818	rBV	1136849	922613	7.85%	1.533%
6	7.034	837	841	844	rBV	2902855	2536344	21.57%	4.215%
7	7.439	906	910	913	rBV	2271946	2125149	18.07%	3.531%
8	8.157	1029	1032	1038	rBV	1272540	1117999	9.51%	1.858%
9	9.233	1211	1215	1218	rBV	3236929	2987805	25.41%	4.965%
10	9.492	1256	1259	1265	rBV	194519	187533	1.59%	0.312%
11	9.622	1277	1281	1288	rVB	332179	313914	2.67%	0.522%
12	9.833	1313	1317	1321	rVB	187966	152829	1.30%	0.254%
13	9.916	1327	1331	1335	rBV	1194942	1050008	8.93%	1.745%
14	10.333	1399	1402	1405	rVB	237644	204639	1.74%	0.340%
15	10.704	1461	1465	1474	rBV	1613418	1519821	12.92%	2.526%
16	10.804	1479	1482	1492	rBV	194394	267522	2.27%	0.445%
17	11.251	1555	1558	1561	rBV	144015	131126	1.12%	0.218%
18	11.404	1580	1584	1587	rBV	1043202	909921	7.74%	1.512%
19	11.786	1645	1649	1653	rBV	5237279	5487895	46.67%	9.120%
20	11.922	1668	1672	1677	rBV2	230314	258932	2.20%	0.430%
21	12.486	1761	1768	1770	rBV	6289523	11759413	100.00%	19.541%
22	12.510	1770	1772	1777	rVV2	6526233	6994297	59.48%	11.623%
23	12.580	1780	1784	1787	rVB	3353454	3049581	25.93%	5.068%
24	12.851	1827	1830	1831	rBV	136287	127822	1.09%	0.212%
25	12.992	1850	1854	1857	rBV	2096631	1981670	16.85%	3.293%
26	13.221	1891	1893	1899	rVB	315164	333615	2.84%	0.554%
27	13.298	1903	1906	1908	rBV	357536	346465	2.95%	0.576%
28	13.321	1908	1910	1915	rVV3	358664	438644	3.73%	0.729%
29	13.369	1915	1918	1922	rVV3	346494	405954	3.45%	0.675%
30	13.410	1923	1925	1930	rVV	189341	191871	1.63%	0.319%
31	13.510	1939	1942	1945	rVB	442712	403441	3.43%	0.670%
32	13.733	1975	1980	1985	rVB2	454916	662325	5.63%	1.101%
33	13.874	2001	2004	2007	rVB	204331	192318	1.64%	0.320%
34	13.904	2007	2009	2014	rVB	141205	146848	1.25%	0.244%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.968	2017	2020	2023	rBV	385651	312967	2.66%	0.520%
36	14.051	2030	2034	2037	rBV	902870	858981	7.30%	1.427%
37	15.545	2284	2288	2295	rVB2	626192	880477	7.49%	1.463%

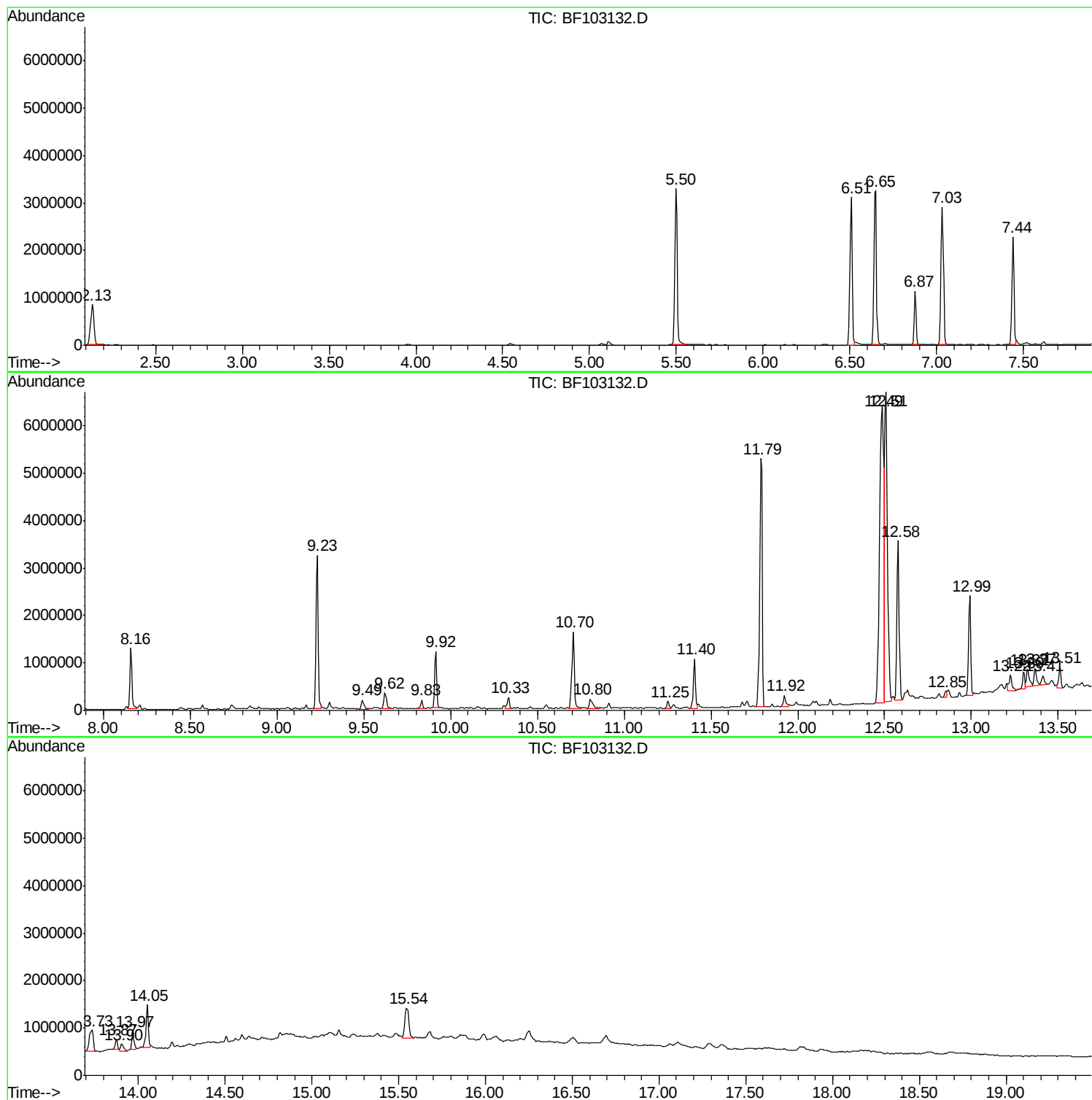
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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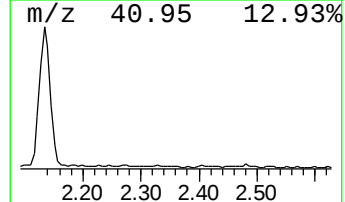
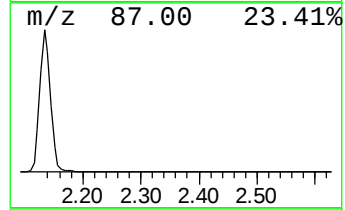
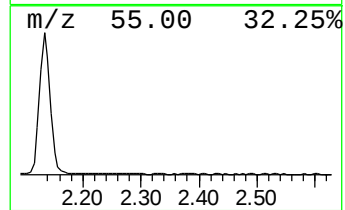
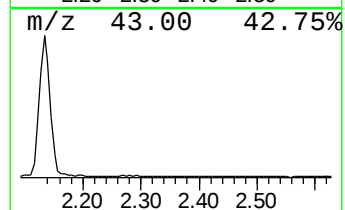
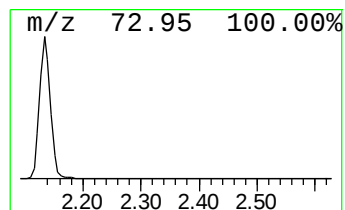
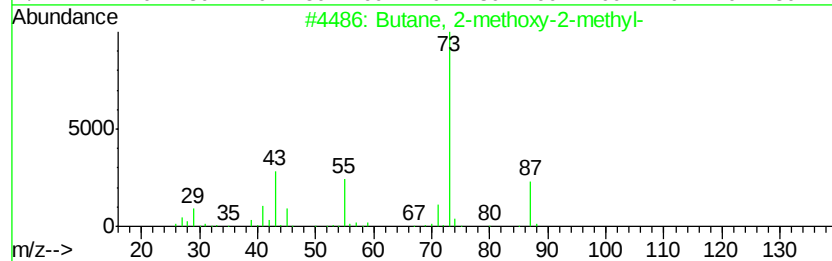
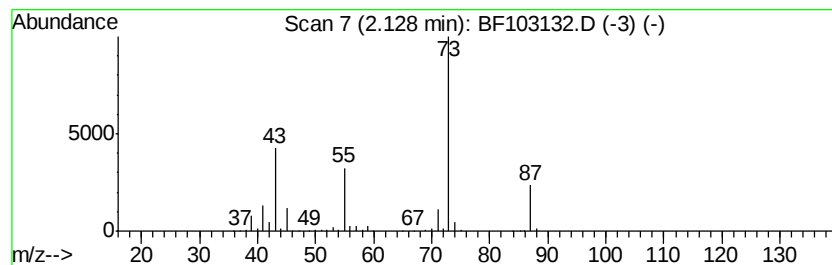
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.13	23.83 ng	1099160	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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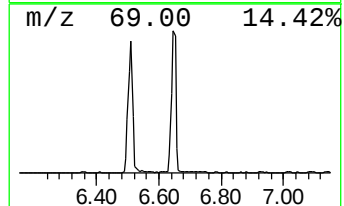
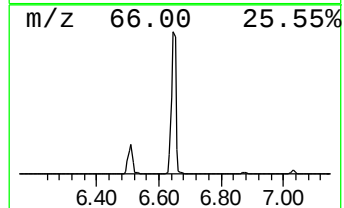
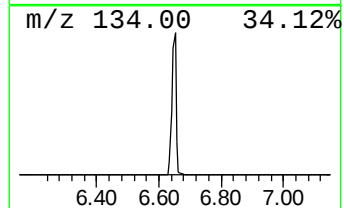
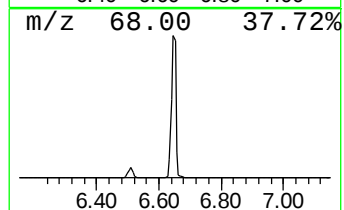
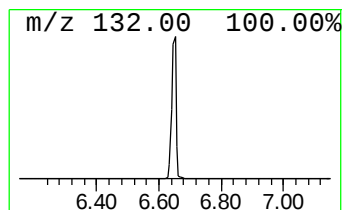
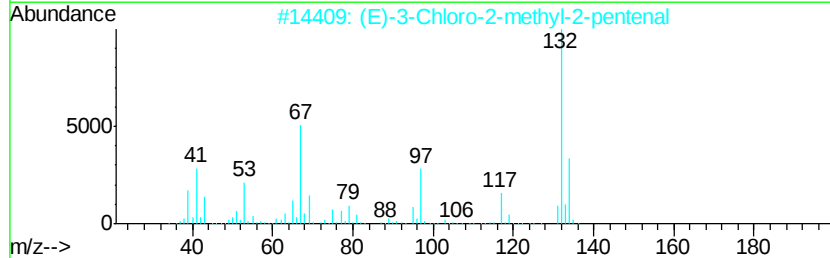
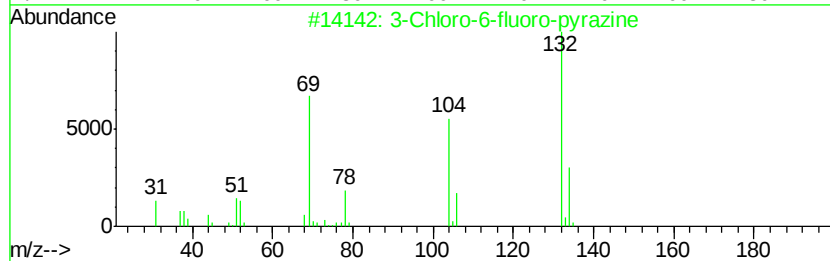
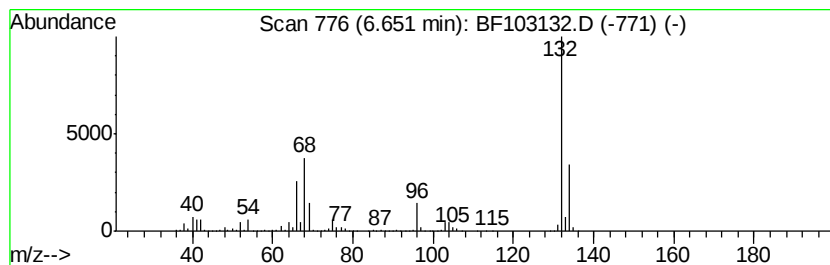
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.65 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	71.27 ng	3287820	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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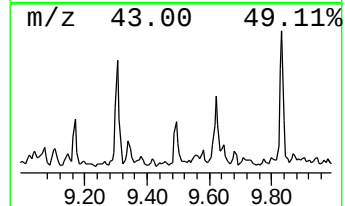
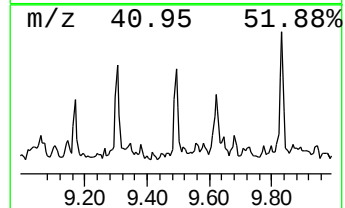
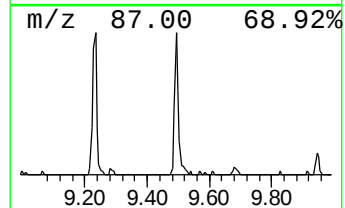
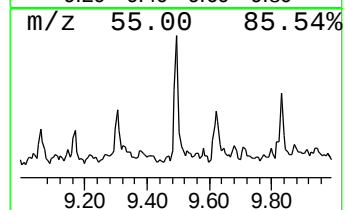
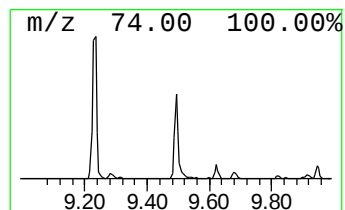
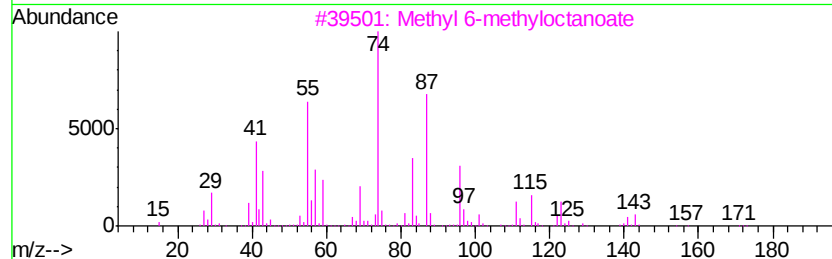
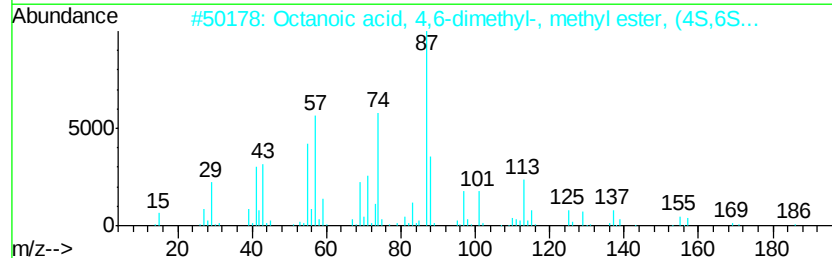
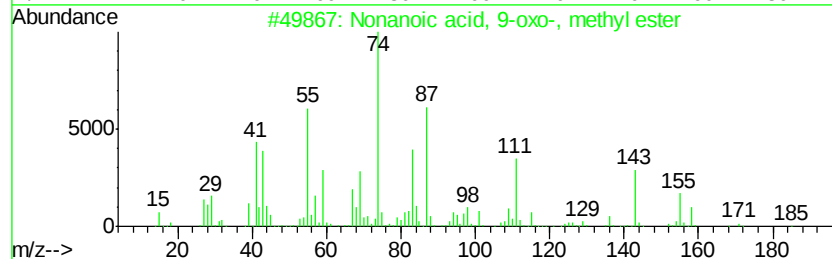
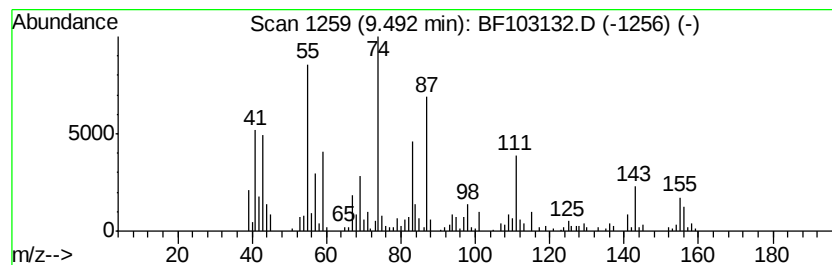
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Nonanoic acid, 9-oxo-, meth... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.49	3.57 ng	187533	Acenaphthene-d10	9.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanoic acid, 9-oxo-, methyl ester	186	C10H18O3	001931-63-1	80		
2	Octanoic acid, 4,6-dimethyl-, me...	186	C11H22O2	002553-96-0	43		
3	Methyl 6-methyloctanoate	172	C10H20O2	005129-62-4	40		
4	Methyl 8-methyl-nonanoate	186	C11H22O2	1000336-43-6	38		
5	Pentanoic acid, 4-methyl-, methy...	130	C7H14O2	002412-80-8	38		



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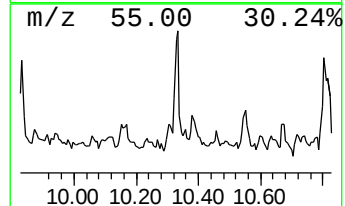
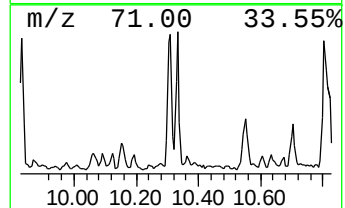
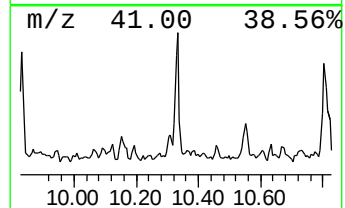
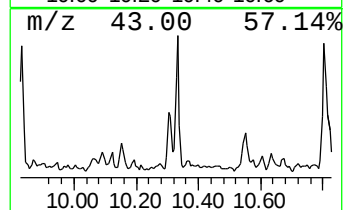
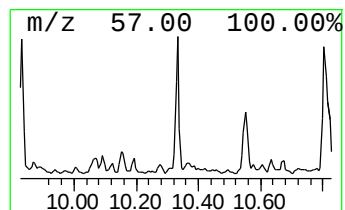
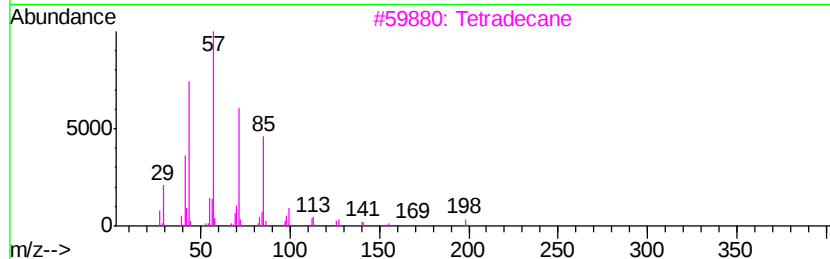
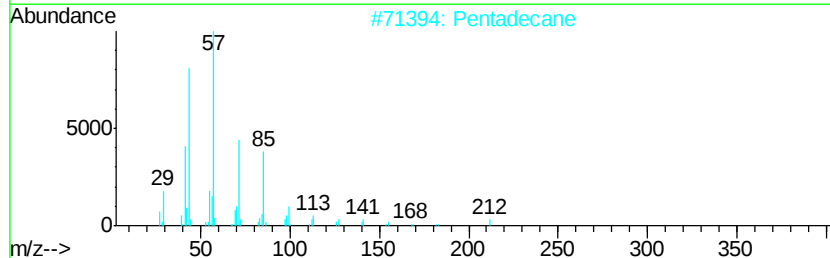
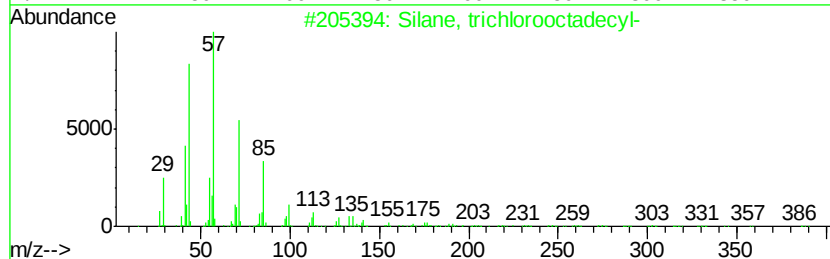
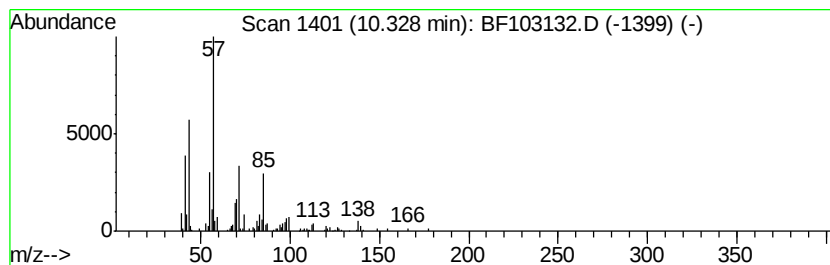
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Silane, trichlorooctadecyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	3.90 ng	204639	Acenaphthene-d10	9.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Silane, trichlorooctadecyl-	386 C18H37Cl3Si	000112-04-9 90
2		Pentadecane	212 C15H32	000629-62-9 90
3		Tetradecane	198 C14H30	000629-59-4 90
4		Hexadecane	226 C16H34	000544-76-3 87
5		Tetracosane	338 C24H50	000646-31-1 83



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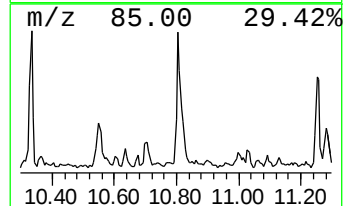
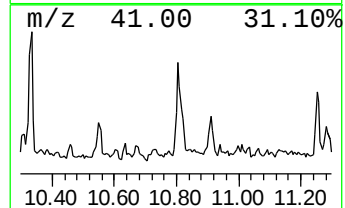
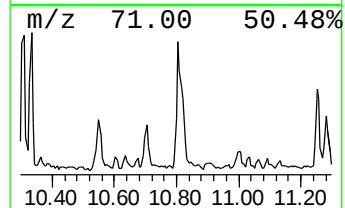
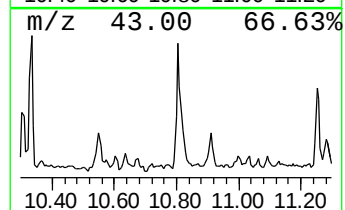
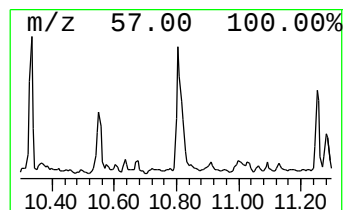
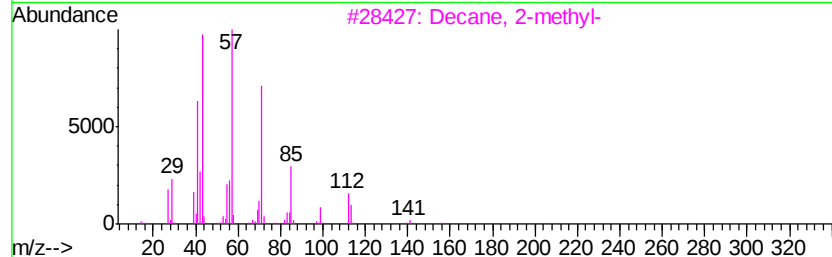
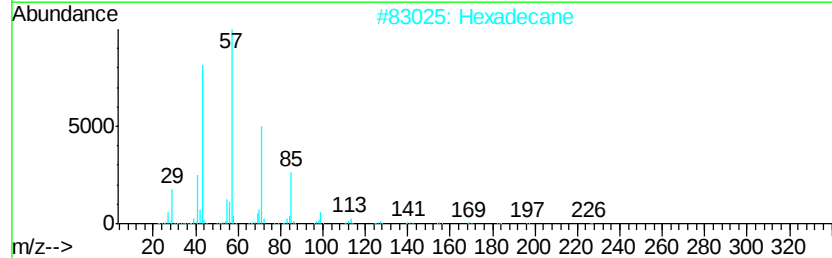
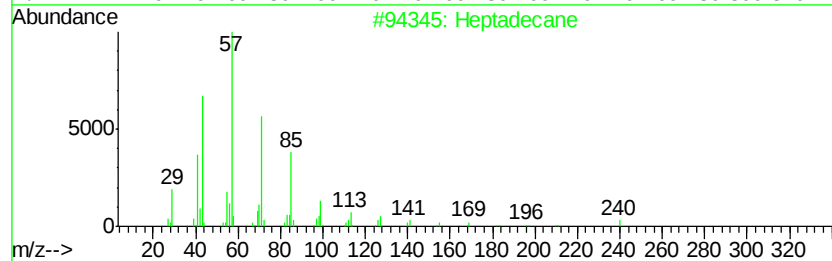
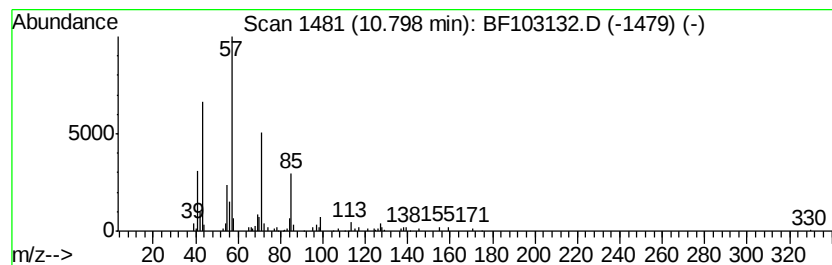
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Heptadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	5.88 ng	267522	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	90
2	Hexadecane		226	C16H34	000544-76-3	90
3	Decane, 2-methyl-		156	C11H24	006975-98-0	74
4	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	72
5	Decane, 3-methyl-		156	C11H24	013151-34-3	68



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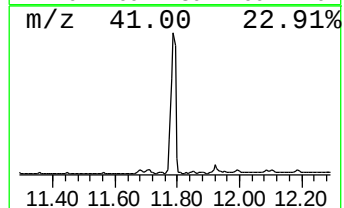
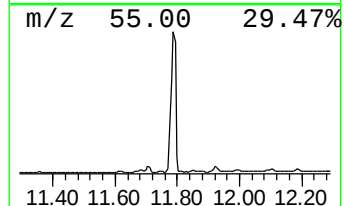
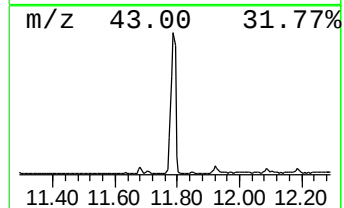
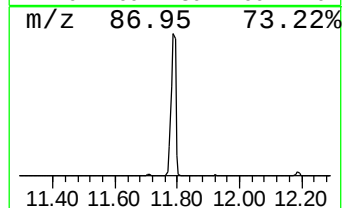
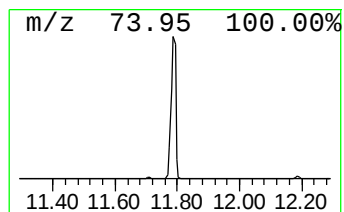
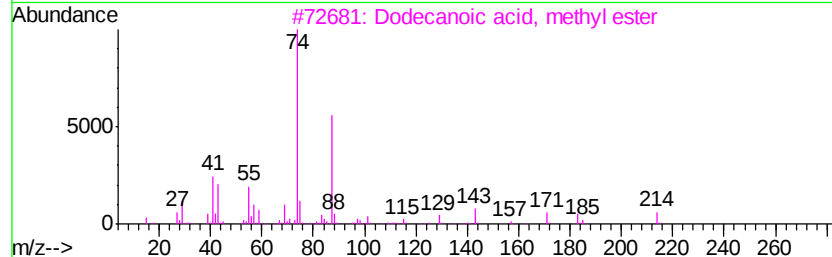
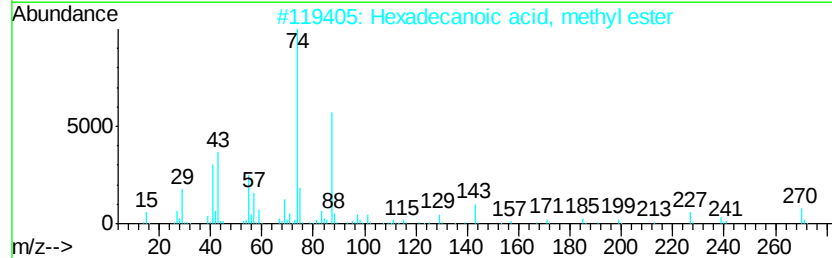
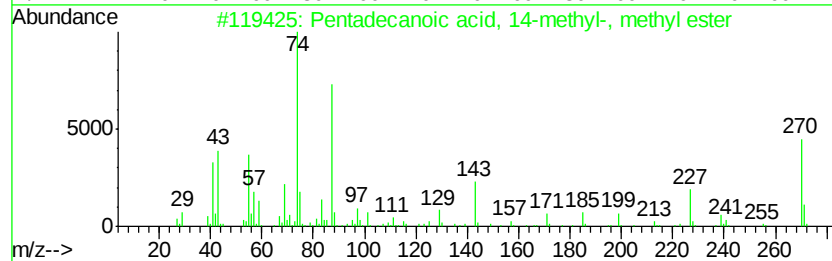
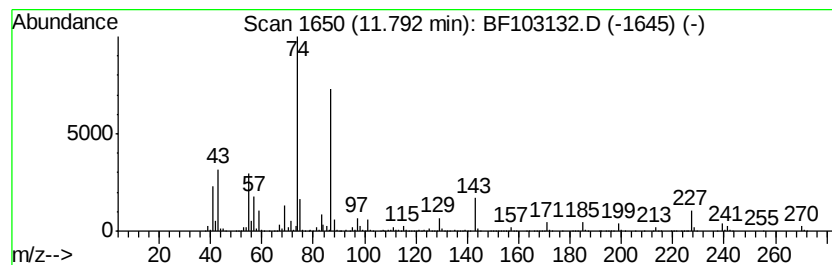
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentadecanoic acid, 14-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	120.62 ng	5487900	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	94
2		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	91
3		Dodecanoic acid, methyl ester	214	C13H26O2	000111-82-0	90
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
5		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022118\
Data File : BF103132.D
Acq On : 21 Feb 2018 13:24
Operator : SJ/JU
Sample : J1404-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-022018

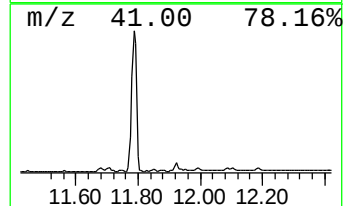
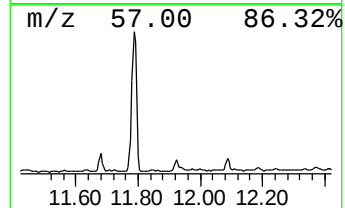
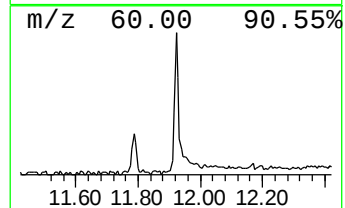
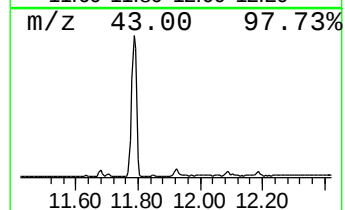
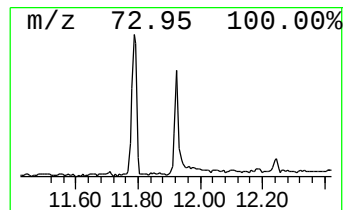
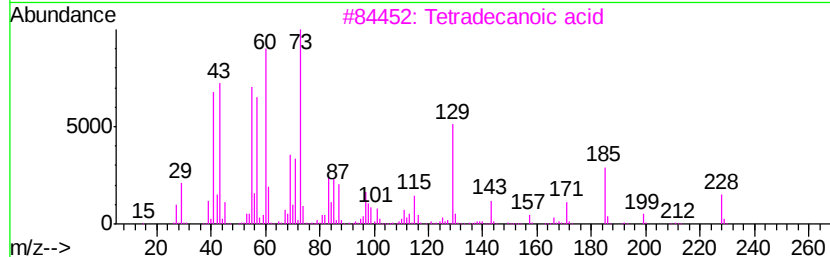
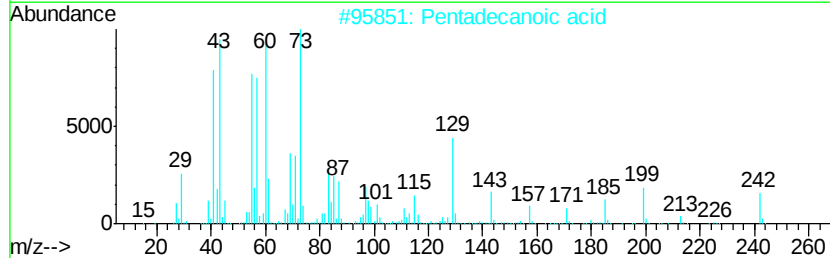
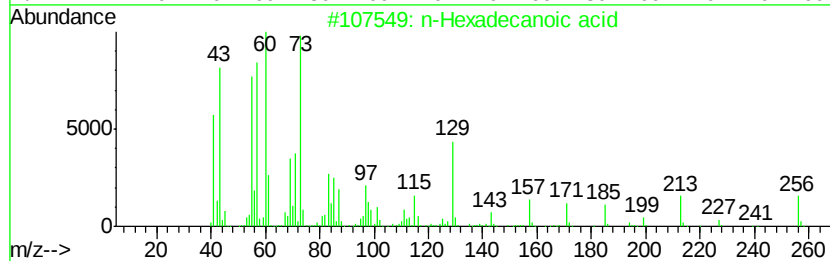
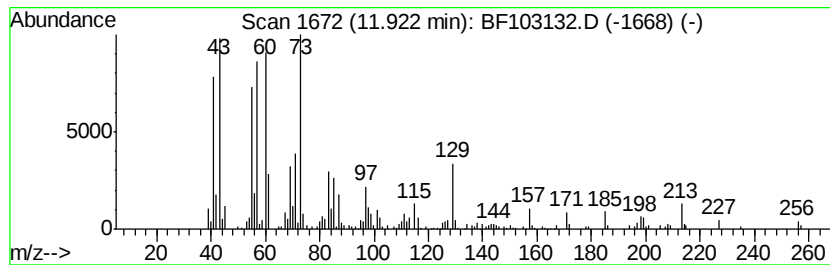
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	5.69 ng	258932	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4		Tridecanoic acid	214	C13H26O2	000638-53-9	72
5		Dodecanoic acid	200	C12H24O2	000143-07-7	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022118\
Data File : BF103132.D
Acq On : 21 Feb 2018 13:24
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Sample : J1404-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

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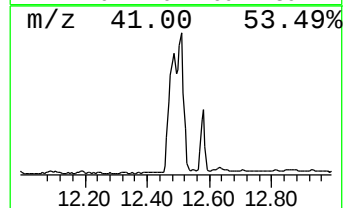
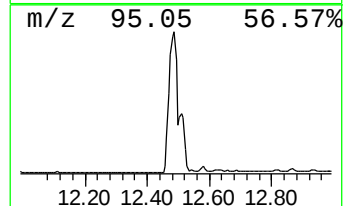
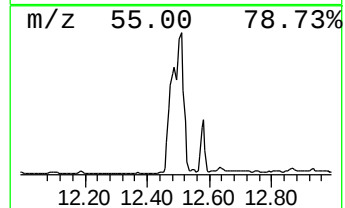
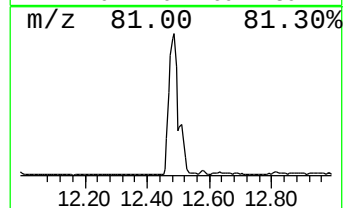
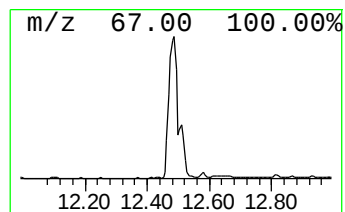
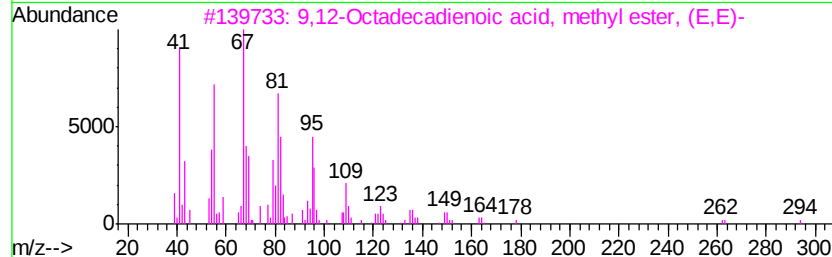
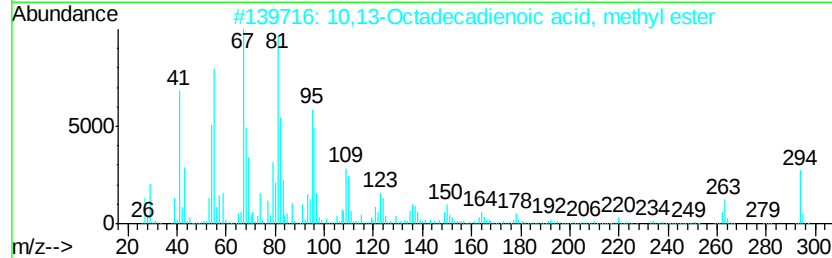
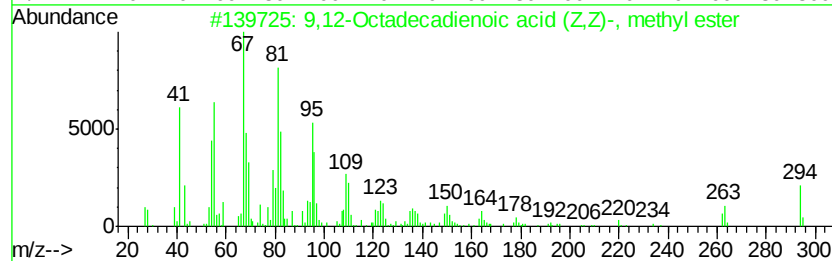
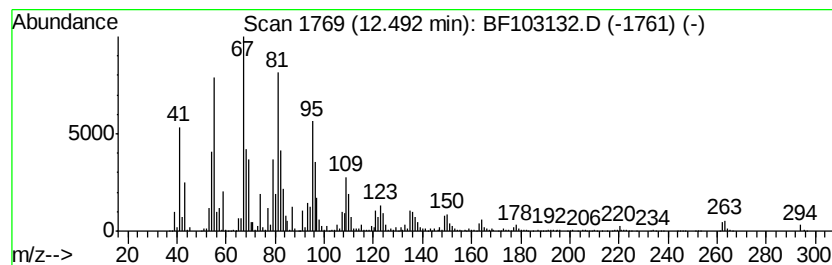
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9,12-Octadecadienoic acid (... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	258.47 ng	11759400	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
2		10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99
3		9,12-Octadecadienoic acid, methv...	294	C19H34O2	002566-97-4	99
4		8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99
5		Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99



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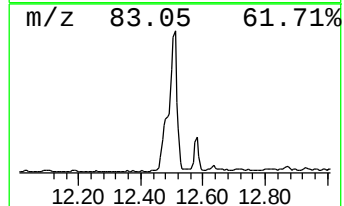
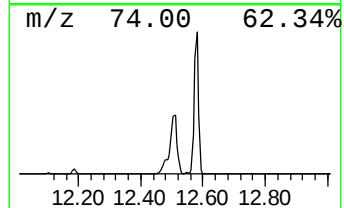
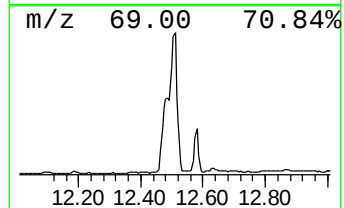
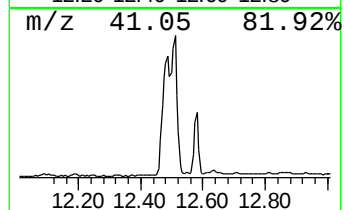
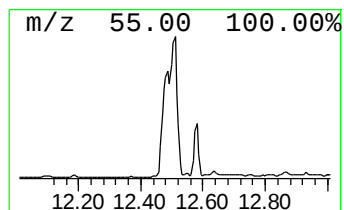
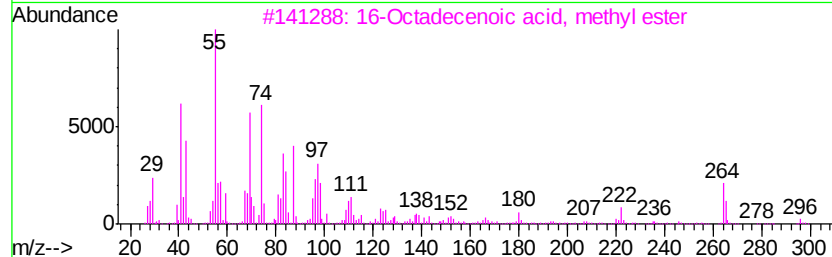
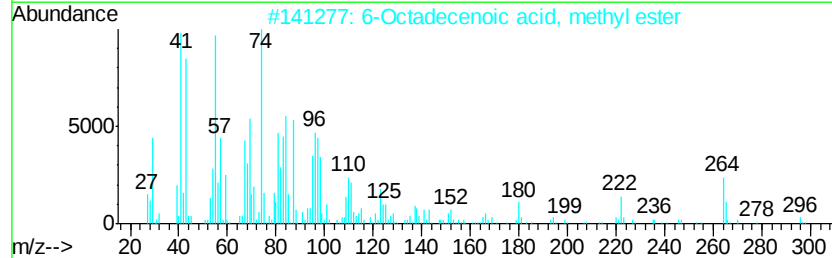
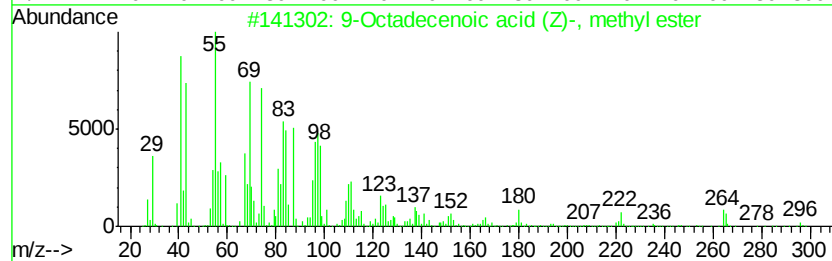
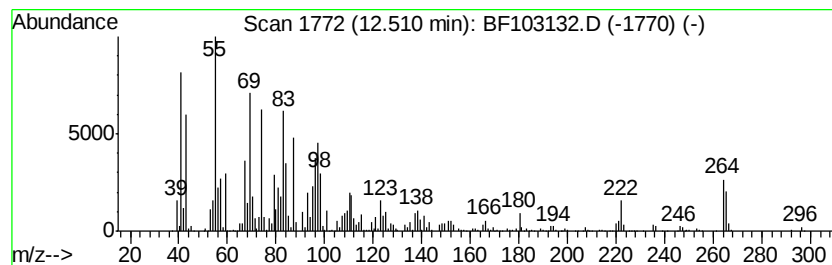
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.51	153.73 ng	6994300	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2		6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	99
3		16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	98
4		13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	98
5		6-Octadecenoic acid, methyl este...	296	C19H36O2	002777-58-4	96



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022118\
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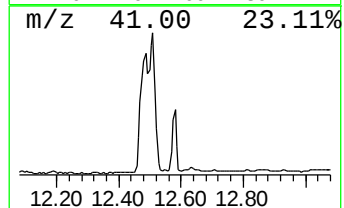
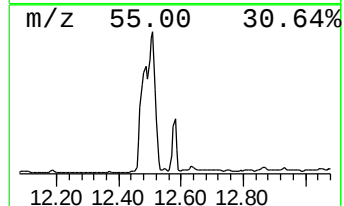
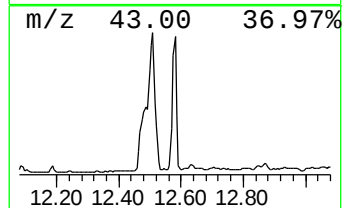
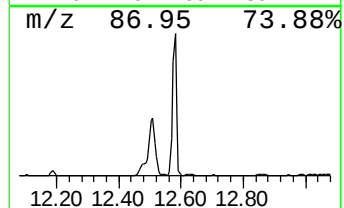
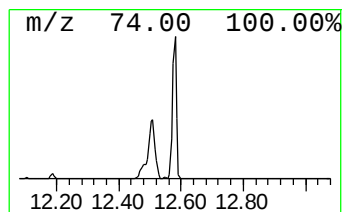
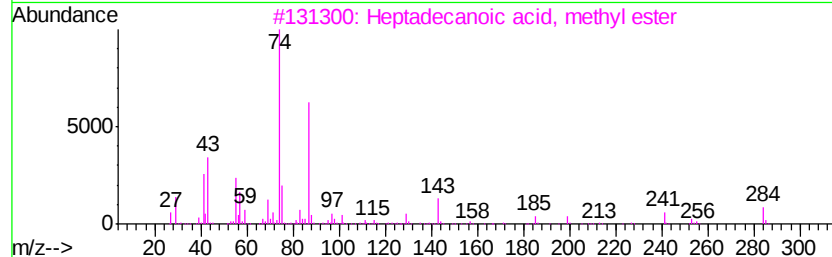
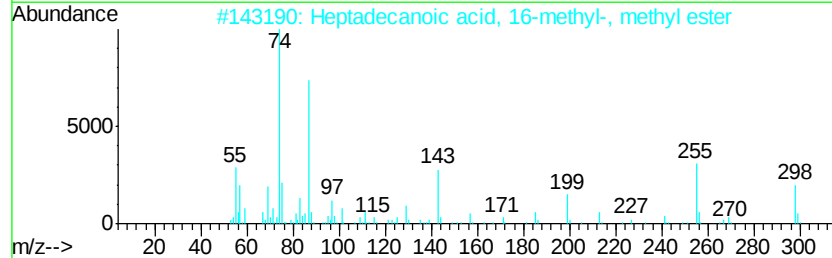
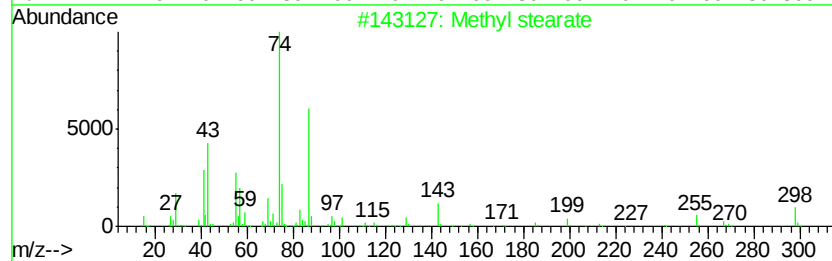
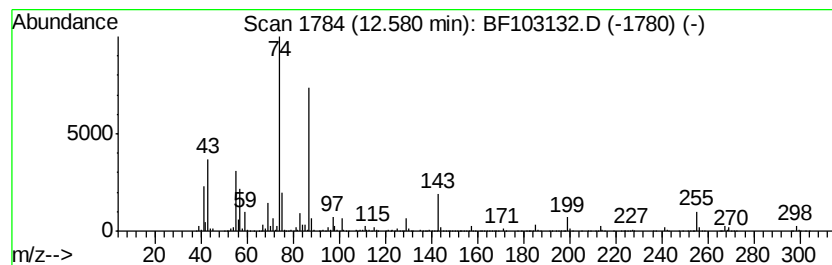
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Methyl stearate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.58	67.03 ng	3049580	Phenanthrene-d10		11.40	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl stearate	298	C19H38O2	000112-61-8	97
2		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	91
3		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	91
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	90
5		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	90



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ALS Vial : 7 Sample Multiplier: 1

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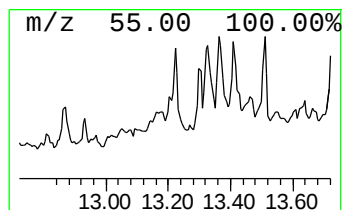
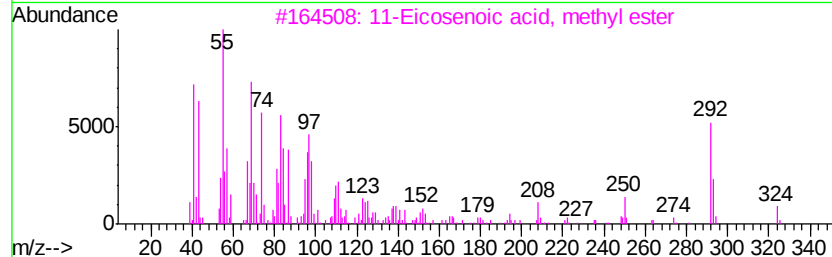
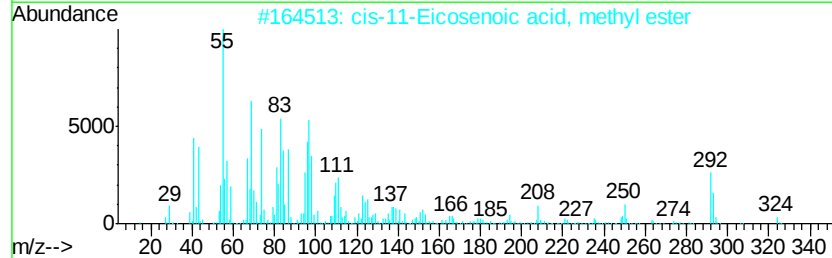
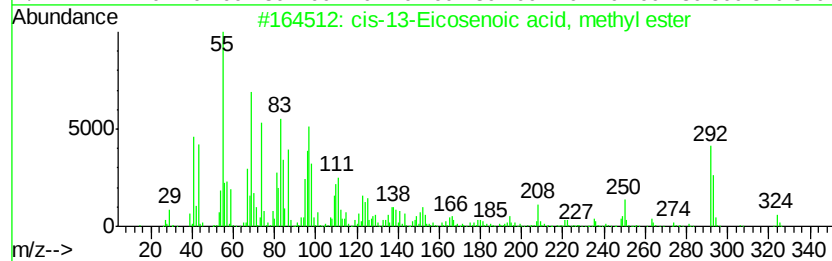
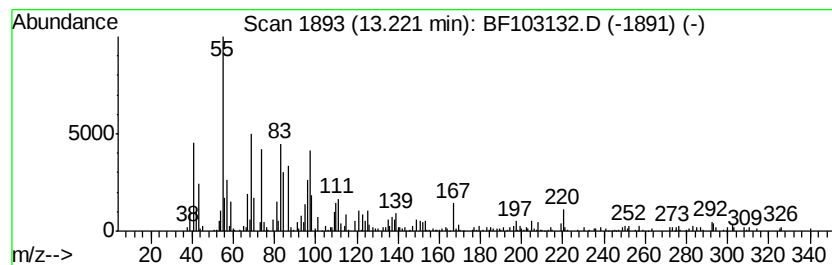
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 cis-13-Eicosenoic acid, met... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	7.77 ng	333615	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-13-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-52-1	91
2		cis-11-Eicosenoic acid, methyl e...	324	C21H40O2	1000333-63-8	90
3		11-Eicosenoic acid, methyl ester	324	C21H40O2	003946-08-5	83
4		cis-10-Heptadecenoic acid, methyl...	282	C18H34O2	1000333-62-1	70
5		Cyclopropaneoctanoic acid, 2-hex...	282	C18H34O2	010152-61-1	62



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022118\
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ALS Vial : 7 Sample Multiplier: 1

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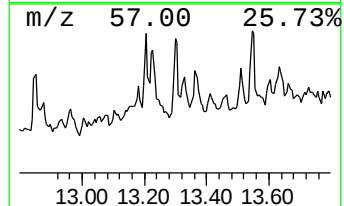
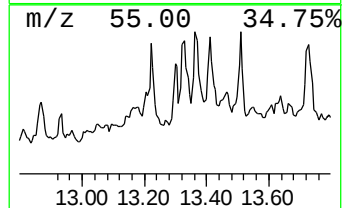
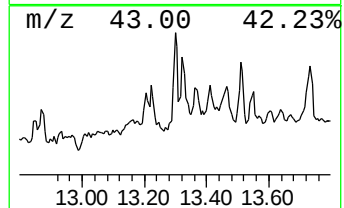
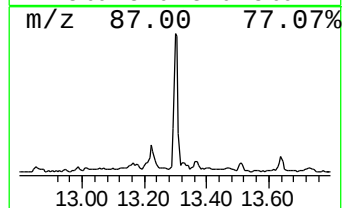
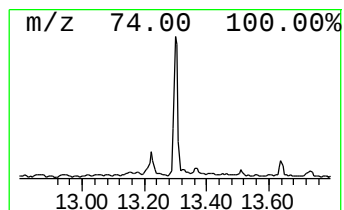
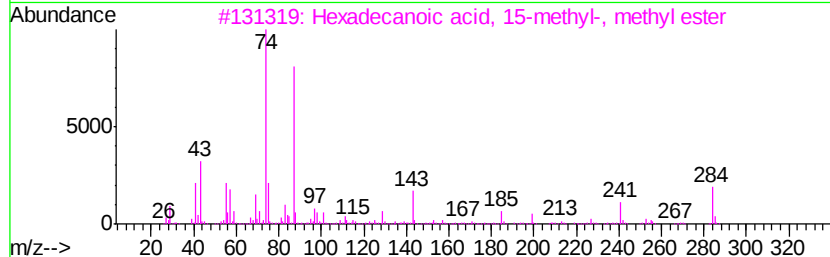
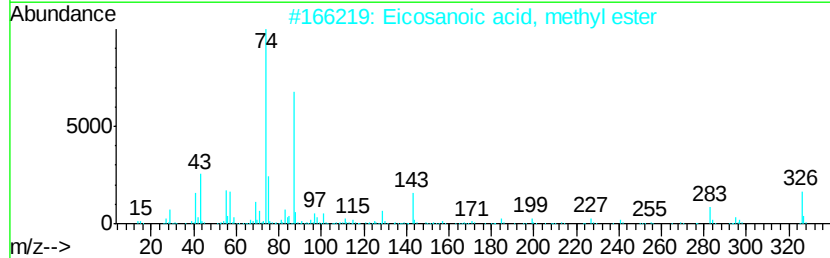
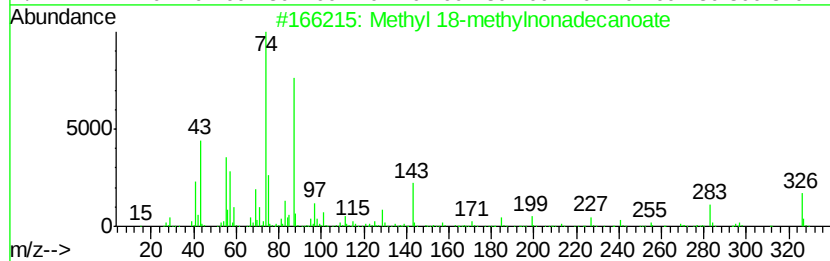
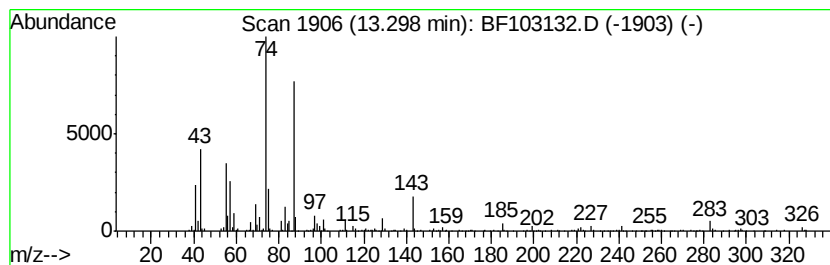
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Methyl 18-methylnonadecanoate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.30	8.07 ng	346465	Chrysene-d12	14.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 18-methylnonadecanoate	326	C21H42O2	1000352-20-6	95
2			Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	95
3			Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	93
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	93
5			Methyl stearate	298	C19H38O2	000112-61-8	90



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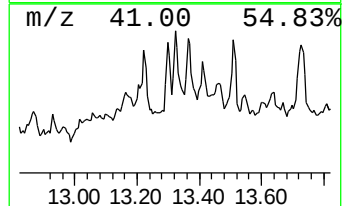
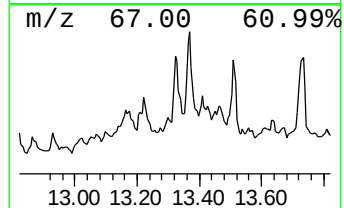
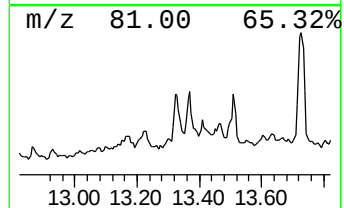
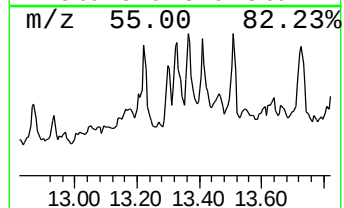
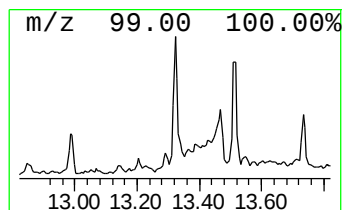
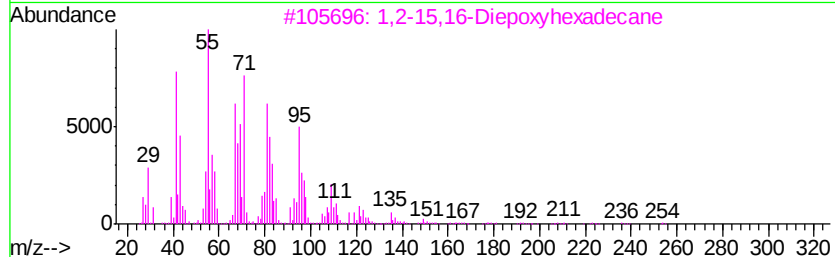
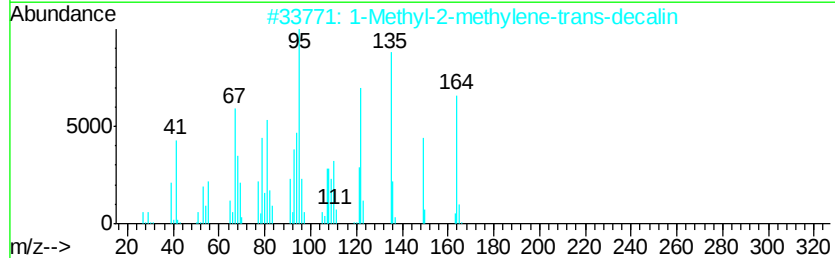
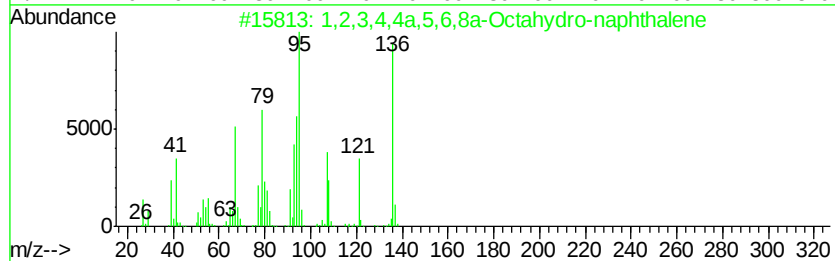
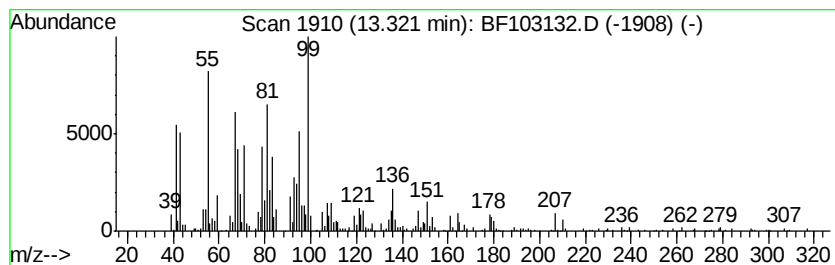
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown13.32 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.32	10.21 ng	438644	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,3,4,4a,5,6,8a-Octahydro-naph...	136	C10H16	031244-58-3	42
2		1-Methyl-2-methylene-trans-decalin	164	C12H20	090548-09-7	42
3		1,2-15,16-Diepoxyhexadecane	254	C16H30O2	1000192-65-0	35
4		3-Methyl-cis-3a,4,7,7a-tetrahydr...	136	C10H16	1000144-04-4	35
5		2H-Pyran-2-one, tetrahydro-6-nonyl-	226	C14H26O2	002721-22-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022118\
Data File : BF103132.D
Acq On : 21 Feb 2018 13:24
Operator : SJ/JU
Sample : J1404-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-022018

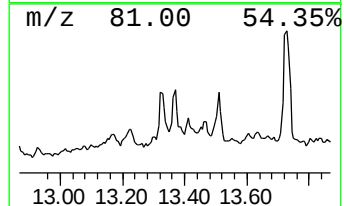
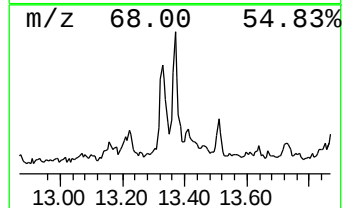
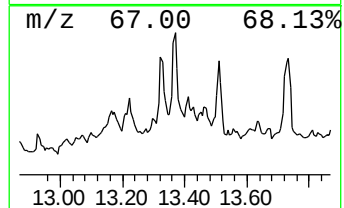
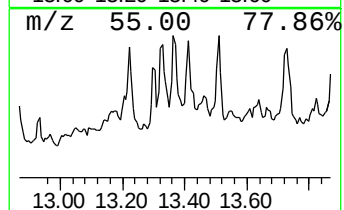
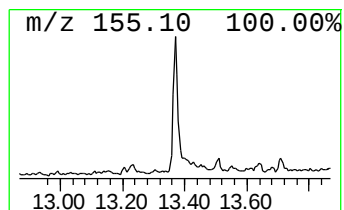
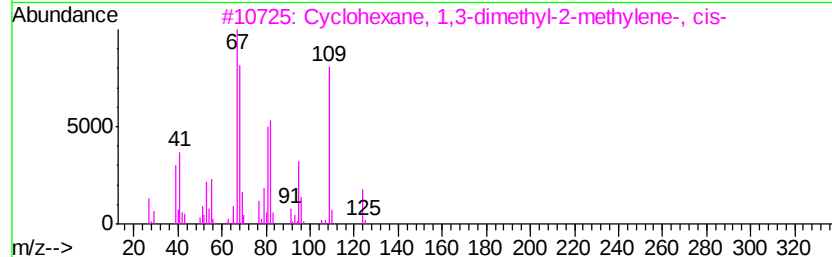
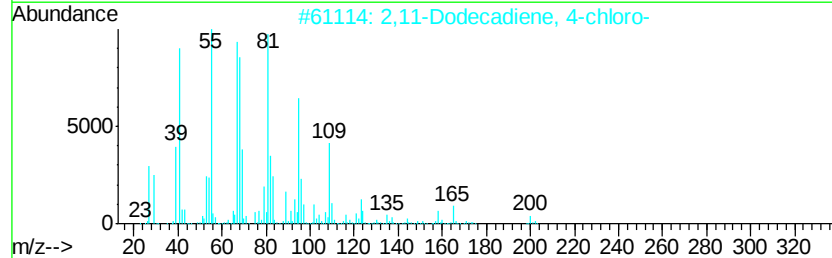
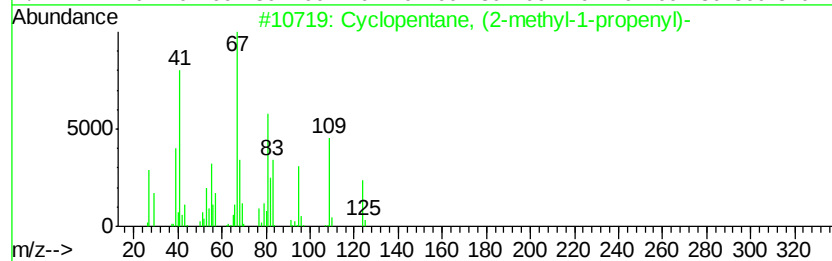
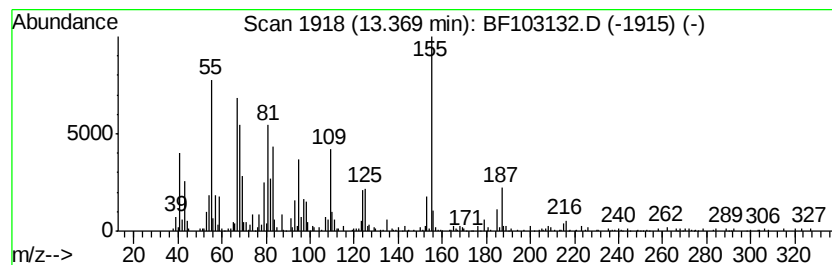
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown13.37 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	9.45 ng	405954	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, (2-methyl-1-propen...	124	C9H16	053366-57-7	30
2		2,11-Dodecadiene, 4-chloro-	200	C12H21Cl	1000155-54-5	30
3		Cyclohexane, 1,3-dimethyl-2-meth...	124	C9H16	019781-47-6	27
4		1-Methylbicyclo[3.2.1]octane	124	C9H16	119972-41-7	27
5		Benzene, 2-fluoro-1-methyl-4-nitro-	155	C7H6FN02	001427-07-2	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022118\
Data File : BF103132.D
Acq On : 21 Feb 2018 13:24
Operator : SJ/JU
Sample : J1404-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-022018

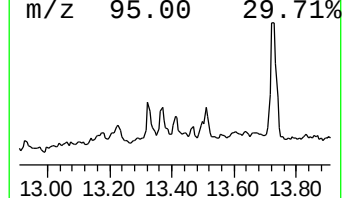
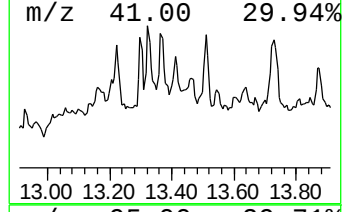
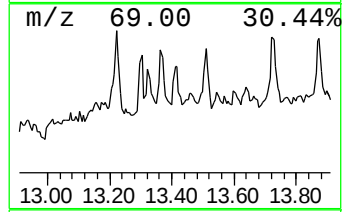
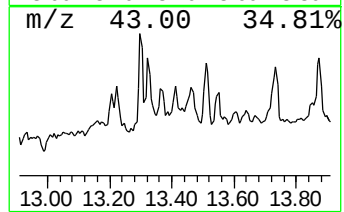
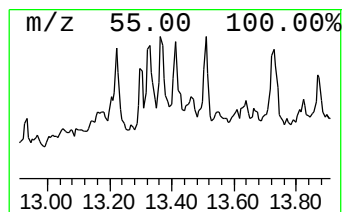
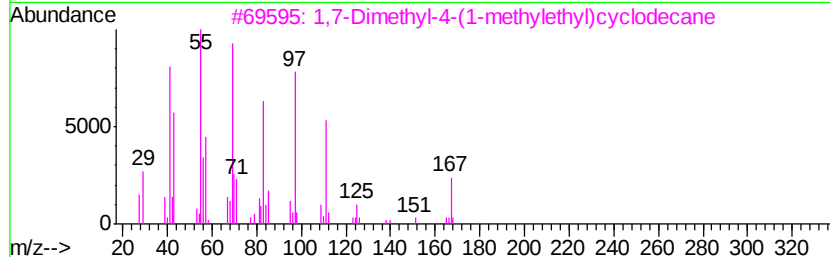
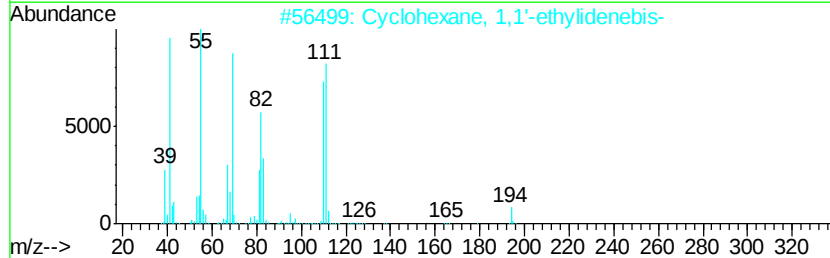
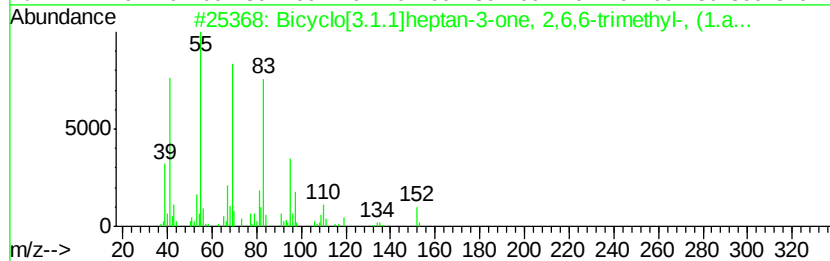
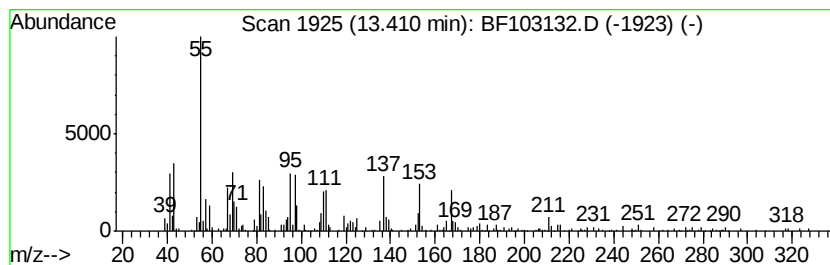
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown13.42 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.41	4.47 ng	191871	Chrysene-d12	14.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	015358-88-0	30
2			Cyclohexane, 1,1'-ethylidenebis-	194	C14H26	002319-61-1	15
3			1,7-Dimethyl-4-(1-methylethyl)cyclodecane	210	C15H30	000645-10-3	14
4			6-Nitroundec-5-ene	199	C11H21NO2	1000192-40-3	14
5			Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	000547-60-4	12



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022118\
Data File : BF103132.D
Acq On : 21 Feb 2018 13:24
Operator : SJ/JU
Sample : J1404-05
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-022018

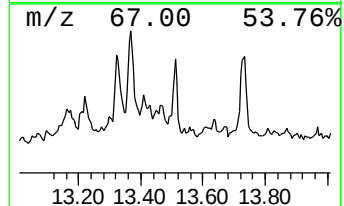
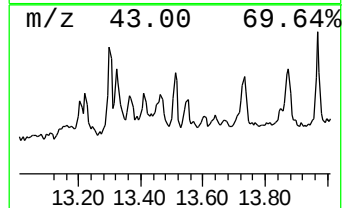
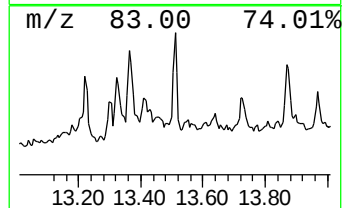
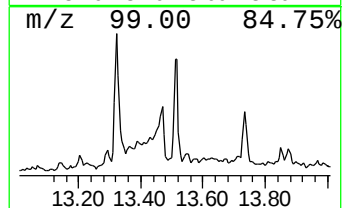
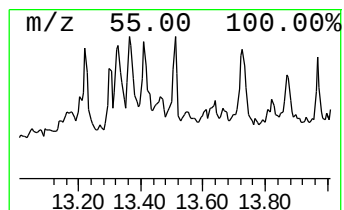
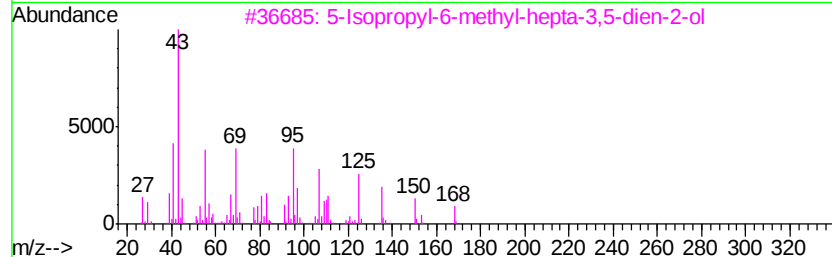
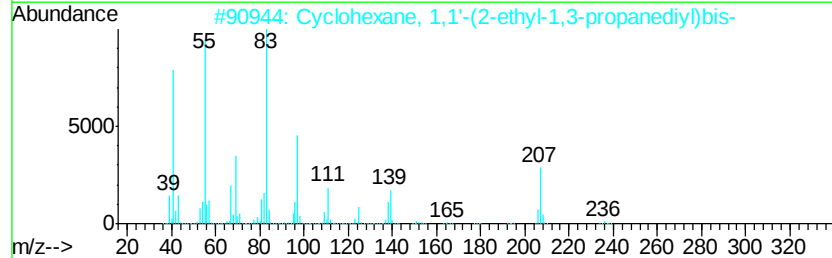
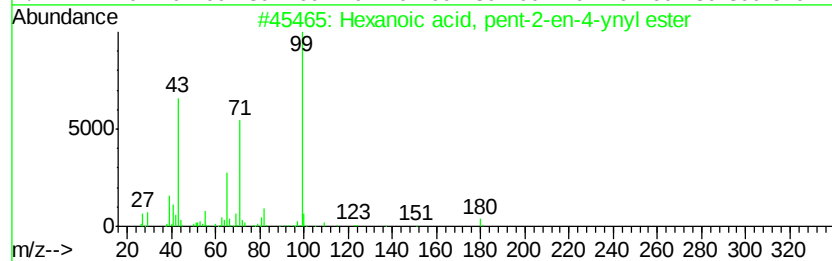
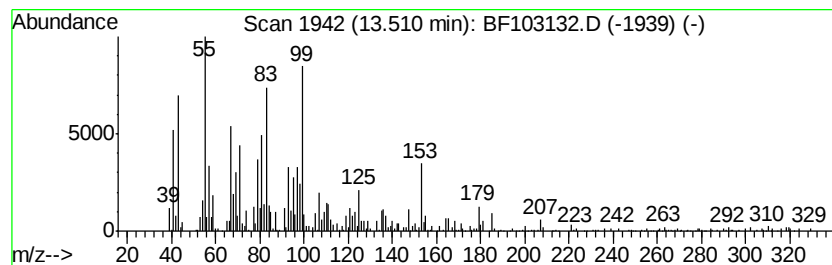
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown13.51 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.51	9.39 ng	403441	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, pent-2-en-4-ynyl ...	180	C11H16O2	1000299-35-3	25
2		Cyclohexane, 1,1'-(2-ethyl-1,3-p...	236	C17H32	054833-34-0	25
3		5-Isopropyl-6-methyl-hepta-3,5-d...	168	C11H20O	1000187-77-8	25
4		Oxirane, 3-[5-(4-azidophenoxy)-3...	287	C16H21N3O2	106929-45-7	25
5		Cycloheptanol, 2-methylene	126	C8H14O	016240-38-3	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022118\
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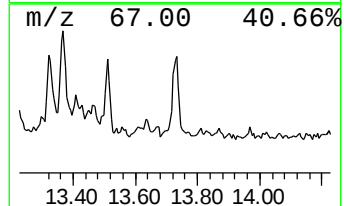
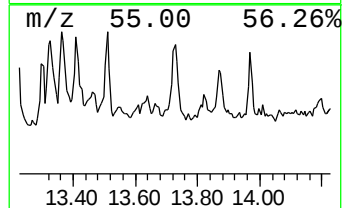
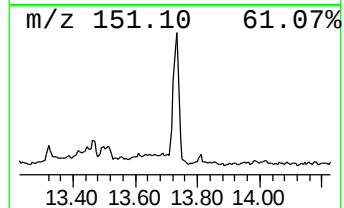
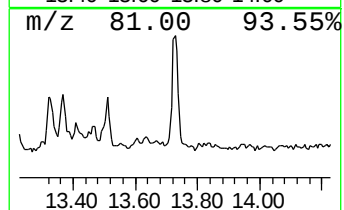
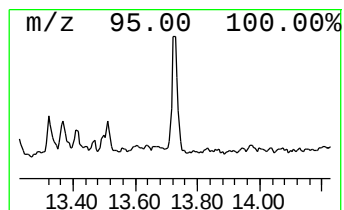
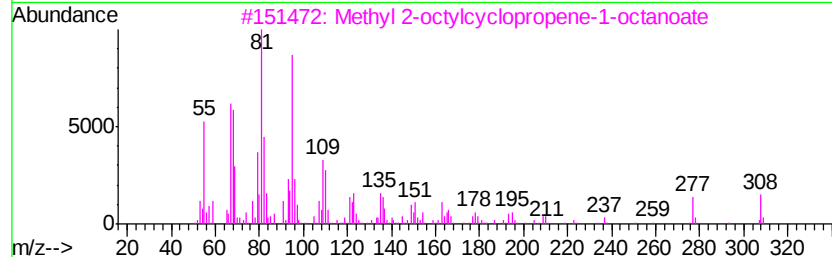
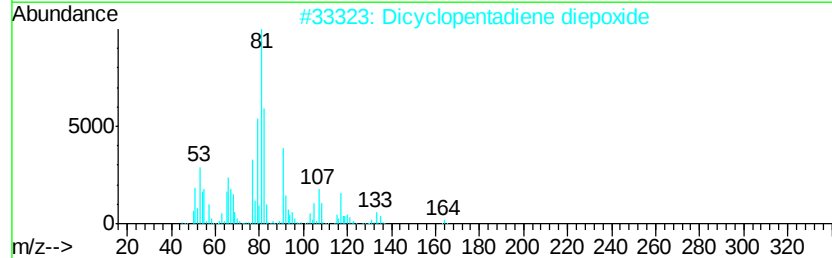
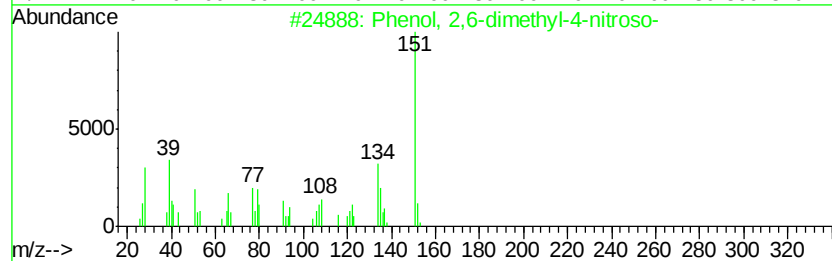
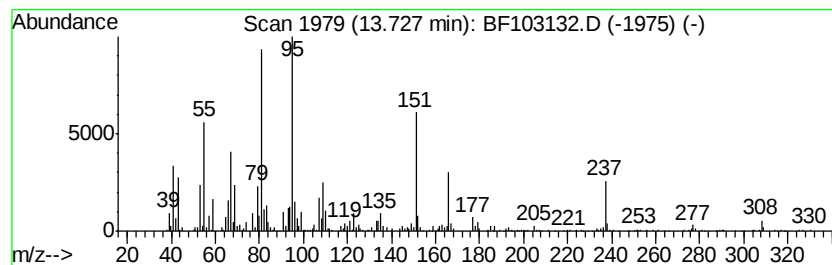
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown13.73 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.73	15.42 ng	662325	Chrysene-d12	14.05		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethyl-4-nitroso-	151	C8H9NO2	013331-93-6	35	
2	Dicyclopentadiene diepoxide	164	C10H12O2	000081-21-0	30	
3	Methyl 2-octylcyclopropene-1-oct...	308	C20H36O2	003220-60-8	20	
4	Guanine	151	C5H5N5O	000073-40-5	11	
5	1,7,7-Trimethylbicyclo[2.2.1]hep...	154	C10H18O	010385-78-1	11	



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022118\
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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ClientSampleId :
TR-05-022018

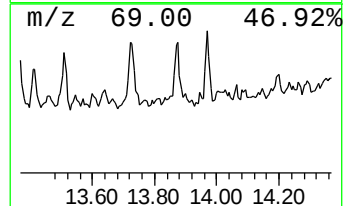
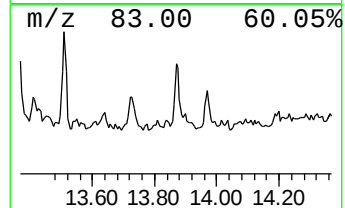
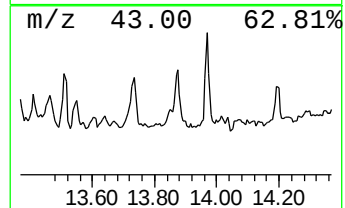
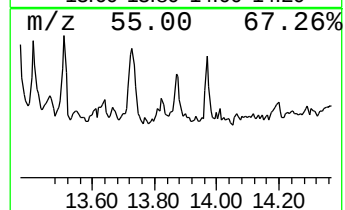
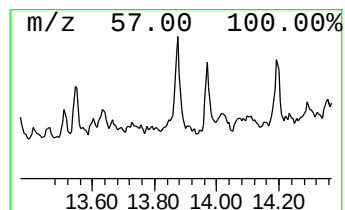
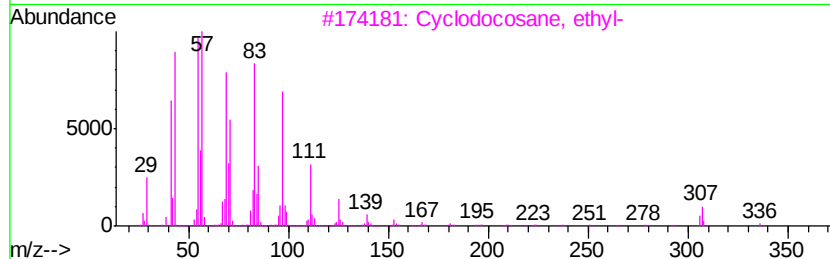
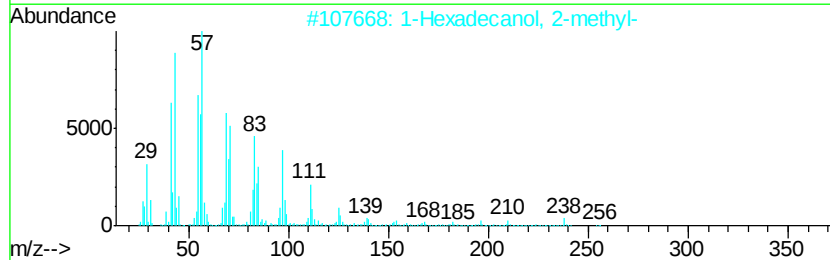
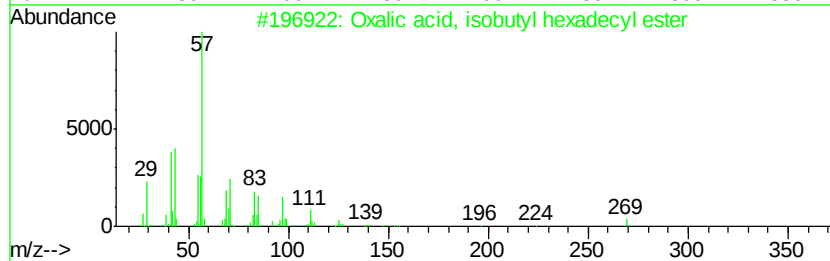
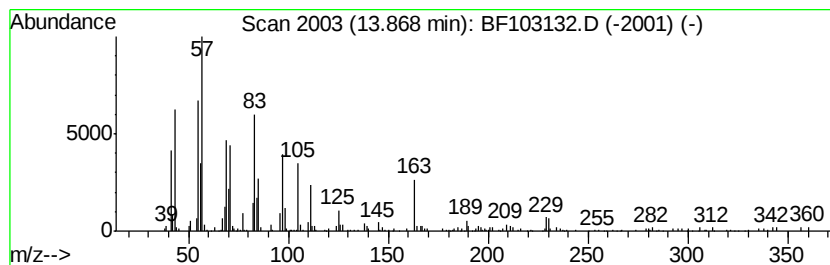
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Oxalic acid, isobutyl hexad... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	4.48 ng	192318	Chrysene-d12	14.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	68
2			1-Hexadecanol, 2-methyl-	256	C17H36O	002490-48-4	58
3			Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	46
4			Oxalic acid, decyl propyl ester	272	C15H28O4	1000309-26-3	43
5			Carbonic acid, isobutyl octadecyl...	370	C23H46O3	1000314-61-5	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022118\
Data File : BF103132.D
Acq On : 21 Feb 2018 13:24
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Misc :
ALS Vial : 7 Sample Multiplier: 1

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ClientSampleId :
TR-05-022018

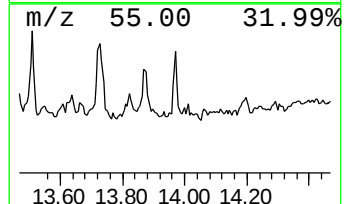
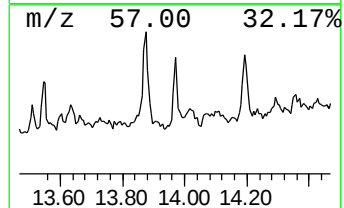
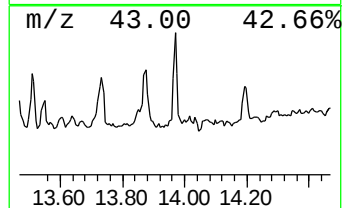
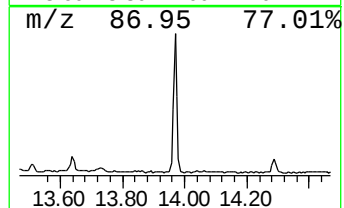
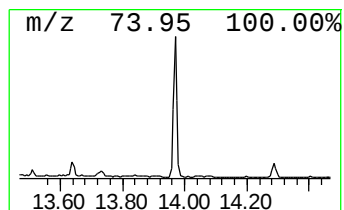
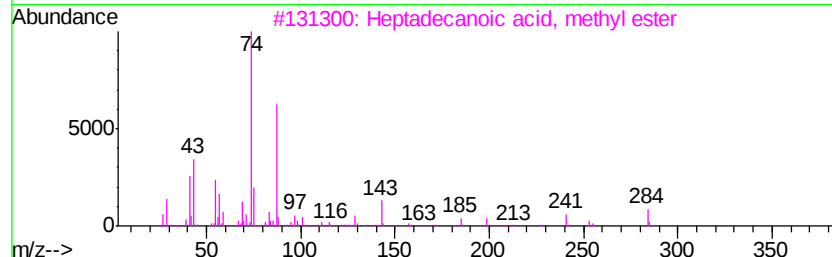
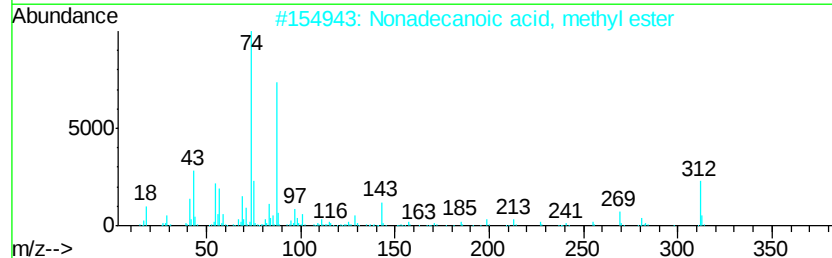
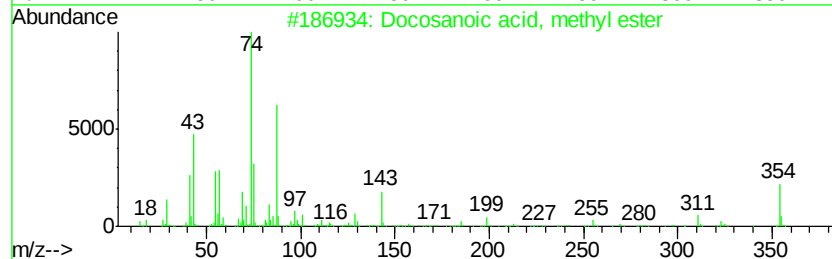
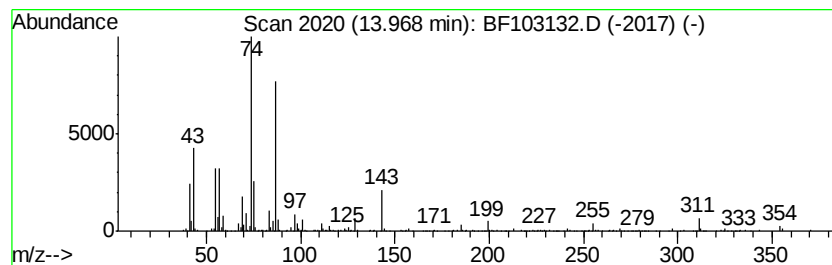
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 Docosanoic acid, methyl ester Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	7.29 ng	312967	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosanoic acid, methyl ester	354	C23H46O2	000929-77-1	96
2		Nonadecanoic acid, methyl ester	312	C20H40O2	001731-94-8	91
3		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	90
4		Methyl stearate	298	C19H38O2	000112-61-8	86
5		Eicosanoic acid, methyl ester	326	C21H42O2	001120-28-1	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022118\
 Data File : BF103132.D
 Acq On : 21 Feb 2018 13:24
 Operator : SJ/JU
 Sample : J1404-05
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-022018

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020318.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
Butane, 2-methoxy...	2.13	23.8	ng	1099160	1	6.87	922613	20.0
unknown6.65	6.65	71.3	ng	3287820	1	6.87	922613	20.0
Nonanoic acid, 9-...	9.49	3.6	ng	187533	3	9.92	1050010	20.0
Silane, trichloro...	10.33	3.9	ng	204639	3	9.92	1050010	20.0
Heptadecane	10.80	5.9	ng	267522	4	11.40	909921	20.0
Pentadecanoic aci...	11.79	120.6	ng	5487900	4	11.40	909921	20.0
n-Hexadecanoic acid	11.92	5.7	ng	258932	4	11.40	909921	20.0
9,12-Octadecadien...	12.49	258.5	ng	11759400	4	11.40	909921	20.0
9-Octadecenoic ac...	12.51	153.7	ng	6994300	4	11.40	909921	20.0
Methyl stearate	12.58	67.0	ng	3049580	4	11.40	909921	20.0
cis-13-Eicosenoic...	13.22	7.8	ng	333615	5	14.05	858981	20.0
Methyl 18-methyln...	13.30	8.1	ng	346465	5	14.05	858981	20.0
unknown13.32	13.32	10.2	ng	438644	5	14.05	858981	20.0
unknown13.37	13.37	9.4	ng	405954	5	14.05	858981	20.0
unknown13.42	13.41	4.5	ng	191871	5	14.05	858981	20.0
unknown13.51	13.51	9.4	ng	403441	5	14.05	858981	20.0
unknown13.73	13.73	15.4	ng	662325	5	14.05	858981	20.0
Oxalic acid, isob...	13.87	4.5	ng	192318	5	14.05	858981	20.0
Docosanoic acid, ...	13.97	7.3	ng	312967	5	14.05	858981	20.0