

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF022420\
 Data File : BF119042.D
 Acq On : 25 Feb 2020 1:25
 Operator : CG/JU
 Sample : L1635-10
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUNR-TS007-01

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.981	489	492	499	rVB	33148	47907	1.04%	0.128%
2	5.387	557	561	583	rBV	3316118	3305924	71.76%	8.851%
3	6.398	728	733	736	rBV	3132285	3214008	69.76%	8.605%
4	6.751	789	793	802	rBV	1212912	1025136	22.25%	2.745%
5	6.910	815	820	823	rBV	3354448	3013917	65.42%	8.069%
6	7.316	884	889	892	rBV	2242174	2216591	48.11%	5.934%
7	8.034	1006	1011	1017	rBV	1543392	1365603	29.64%	3.656%
8	9.110	1188	1194	1197	rBV	4752745	4606909	100.00%	12.334%
9	9.504	1257	1261	1265	rBV	287498	260108	5.65%	0.696%
10	9.704	1286	1295	1303	rBV3	33216	91166	1.98%	0.244%
11	9.786	1304	1309	1313	rBV	1904014	1665095	36.14%	4.458%
12	10.575	1438	1443	1453	rBV	3261202	3039548	65.98%	8.138%
13	11.269	1557	1561	1565	rBV	1900899	1627120	35.32%	4.356%
14	11.798	1648	1651	1664	rVB	484402	517069	11.22%	1.384%
15	11.963	1675	1679	1692	rVB5	57455	147086	3.19%	0.394%
16	12.051	1692	1694	1705	rVB7	42210	76989	1.67%	0.206%
17	12.210	1717	1721	1725	rVB	157042	136534	2.96%	0.366%
18	12.522	1762	1774	1777	rBV	363735	416347	9.04%	1.115%
19	12.580	1780	1784	1792	rBV2	148102	172005	3.73%	0.461%
20	12.851	1825	1830	1834	rBV	4377392	4220771	91.62%	11.300%
21	13.063	1862	1866	1868	rBV	54396	62104	1.35%	0.166%
22	13.698	1969	1974	1979	rBV	195082	240953	5.23%	0.645%
23	13.745	1979	1982	1989	rVB	671499	722459	15.68%	1.934%
24	13.904	2005	2009	2016	rVB	1251769	1147712	24.91%	3.073%
25	14.069	2034	2037	2044	rBV	79005	101996	2.21%	0.273%
26	14.380	2086	2090	2101	rBV	677614	759534	16.49%	2.033%
27	14.669	2137	2139	2148	rVB3	61809	94722	2.06%	0.254%
28	14.739	2148	2151	2155	rVB	61261	60503	1.31%	0.162%
29	14.810	2159	2163	2175	rVB10	56132	116591	2.53%	0.312%
30	14.992	2190	2194	2196	rBV	148250	188087	4.08%	0.504%
31	15.016	2196	2198	2202	rVB2	88209	106079	2.30%	0.284%
32	15.333	2246	2252	2258	rBV	815718	1108279	24.06%	2.967%
33	15.539	2283	2287	2293	rVB2	51825	70693	1.53%	0.189%
34	15.757	2320	2324	2328	rBV2	72968	102848	2.23%	0.275%

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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	15.810	2328	2333	2341	rVB	98680	170978	3.71%	0.458%
36	16.504	2448	2451	2459	rVB7	41026	66762	1.45%	0.179%
37	16.874	2510	2514	2523	rVB8	61774	133649	2.90%	0.358%
38	17.262	2575	2580	2587	rBV7	77393	183204	3.98%	0.490%
39	17.339	2588	2593	2599	rVB9	53583	105548	2.29%	0.283%
40	17.580	2629	2634	2640	rBV10	34432	80759	1.75%	0.216%
41	17.698	2648	2654	2661	rBV4	70414	194245	4.22%	0.520%
42	18.027	2703	2710	2723	rVB4	72375	252399	5.48%	0.676%
43	18.239	2740	2746	2753	rVB7	49246	115415	2.51%	0.309%

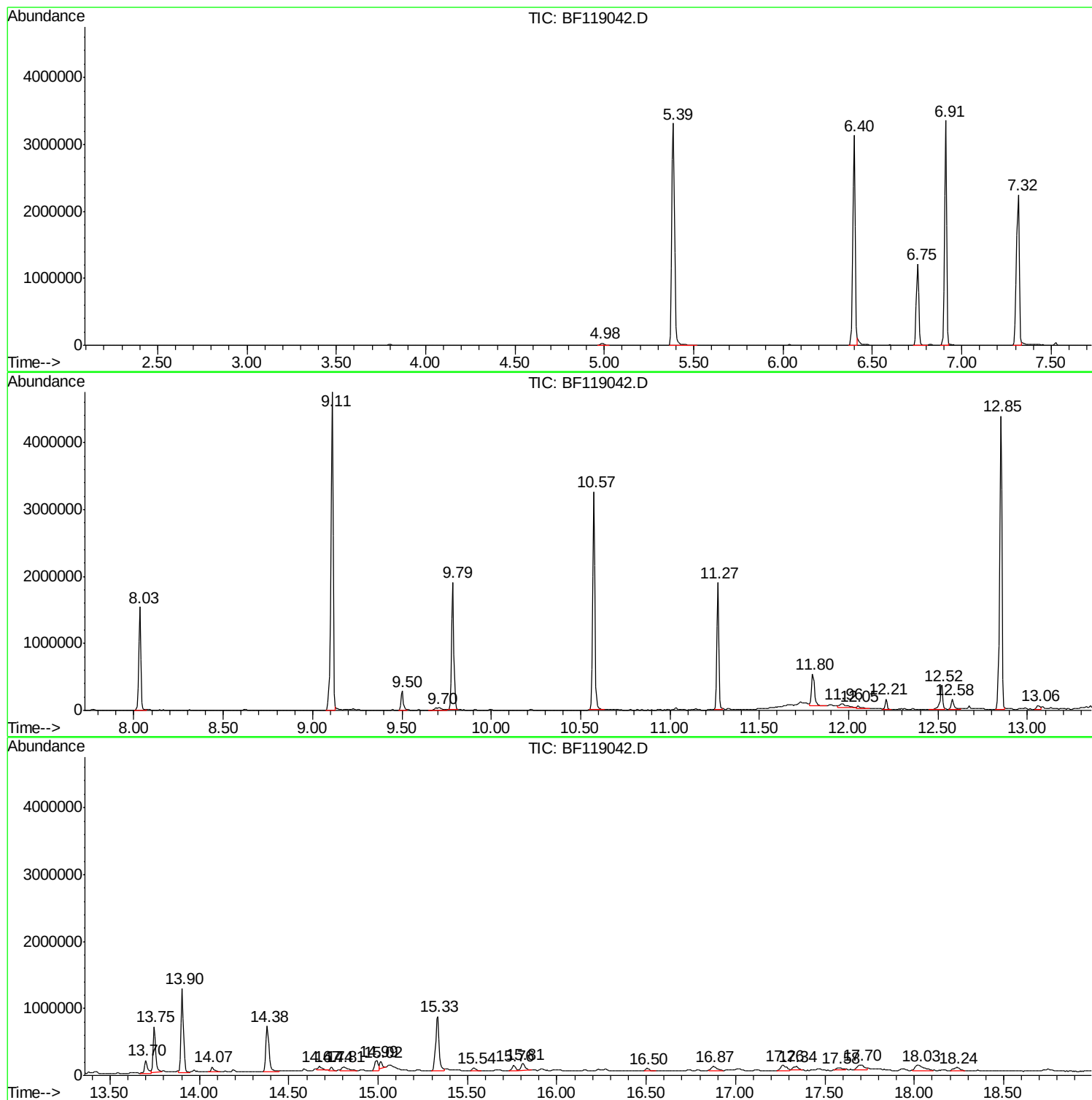
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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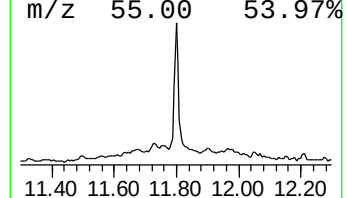
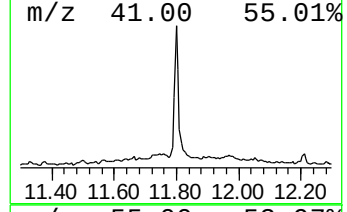
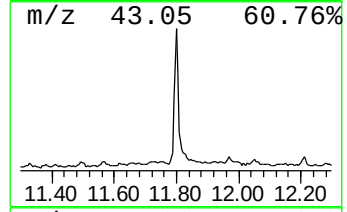
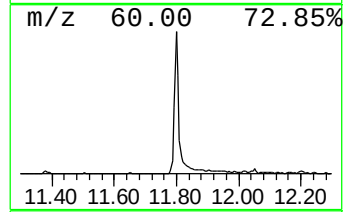
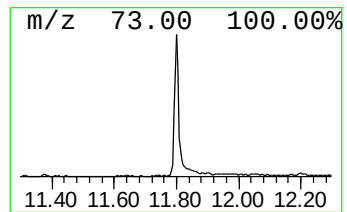
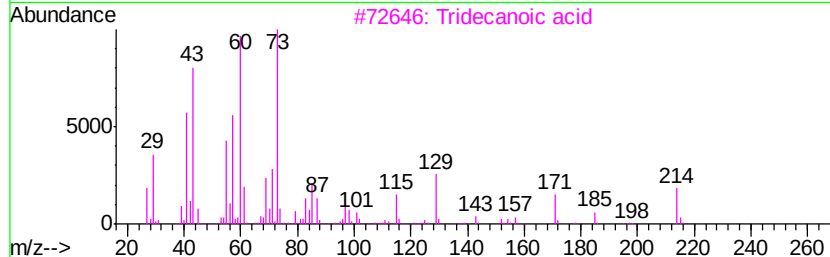
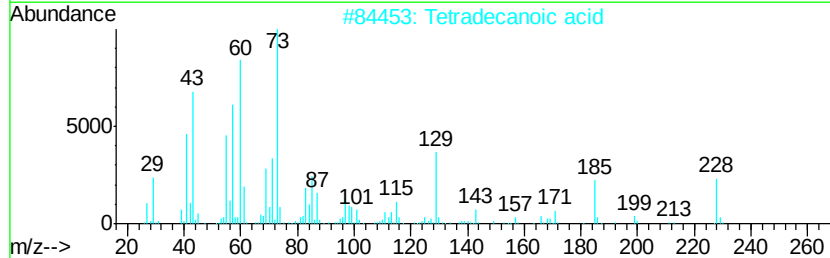
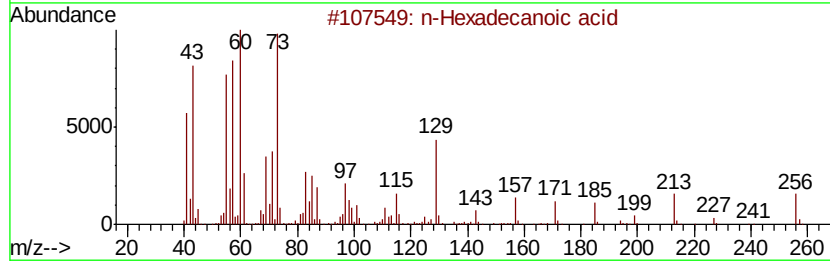
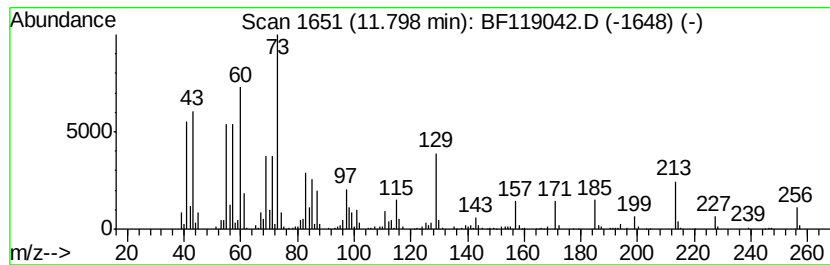
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.80	6.36 ng	517069	Phenanthrene-d10	11.27	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3	Tridecanoic acid	214	C13H26O2	000638-53-9	89
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	62
5	n-Decanoic acid	172	C10H20O2	000334-48-5	60



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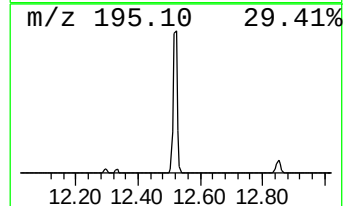
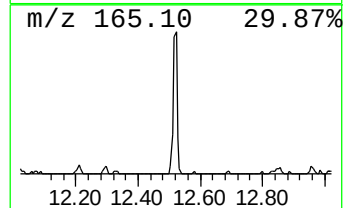
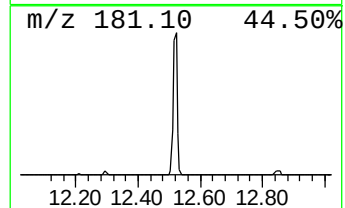
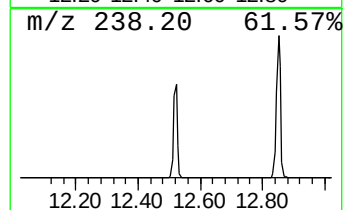
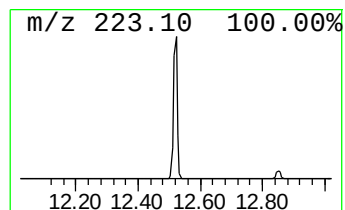
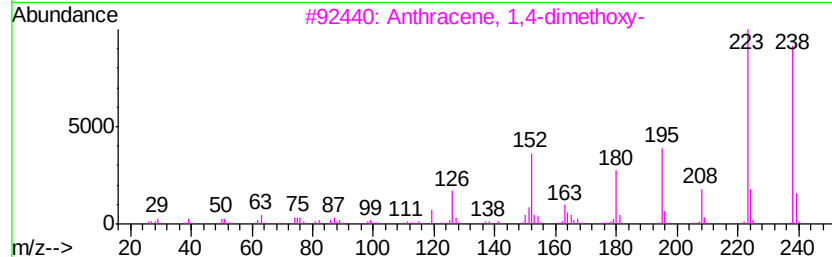
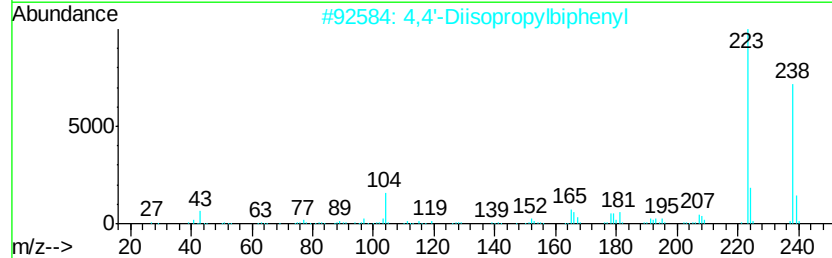
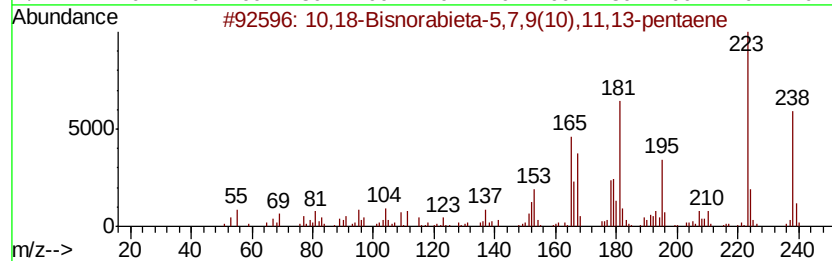
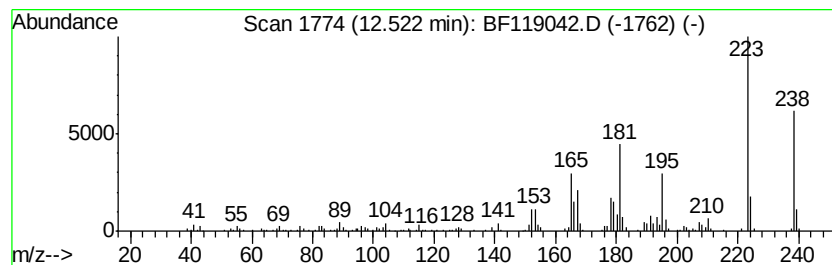
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TIC Integration Parameters: LSCINT.P

Peak Number 3 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.		
12.52	5.12 ng	416347	Phenanthrene-d10		11.27		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99		
2	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	60		
3	Anthracene, 1,4-dimethoxy-	238	C16H14O2	013076-29-4	58		
4	4,4'-Diacetyl biphenyl	238	C16H14O2	000787-69-9	43		
5	9,10-Anthracenedione, 1-amino-	223	C14H9NO2	000082-45-1	43		



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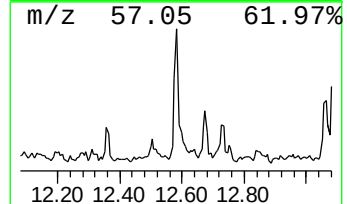
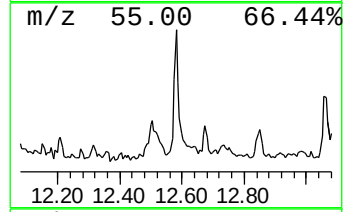
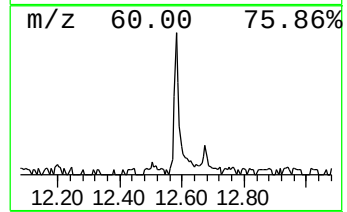
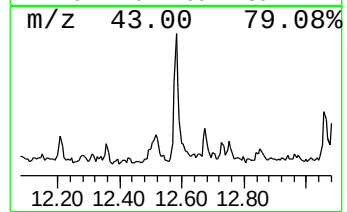
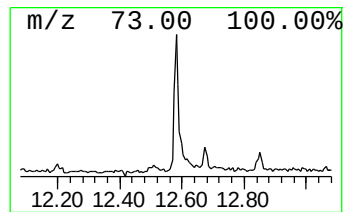
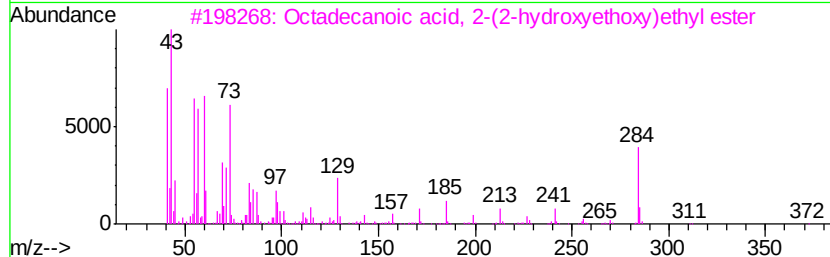
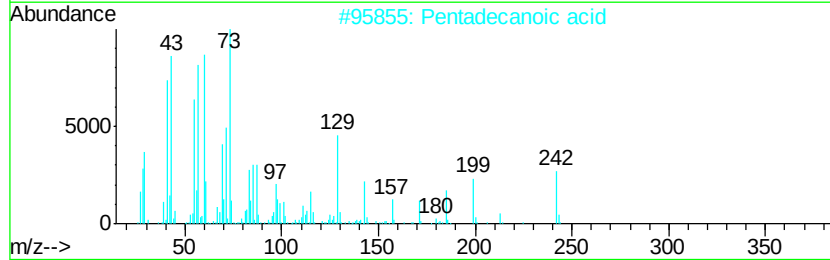
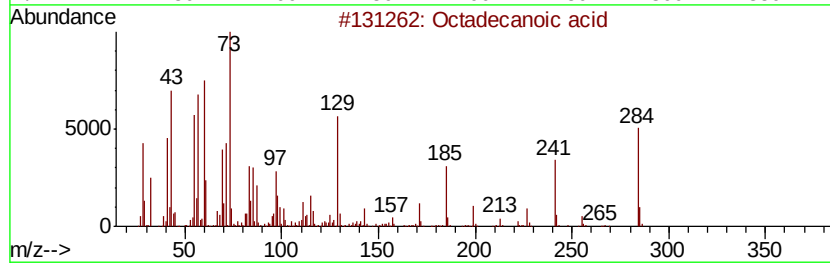
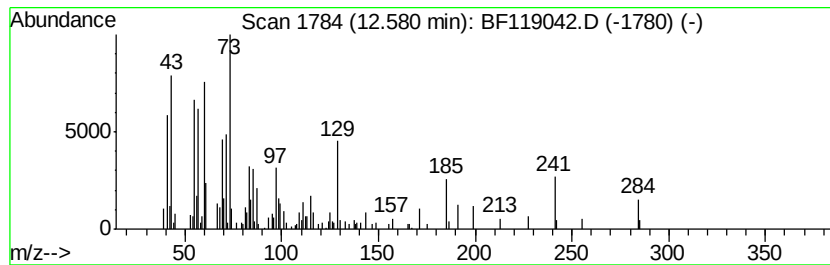
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.		
12.58	2.11 ng	172005	Phenanthrene-d10		11.27		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	89
3			Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	80
4			Undecanoic acid	186	C11H22O2	000112-37-8	74
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	72



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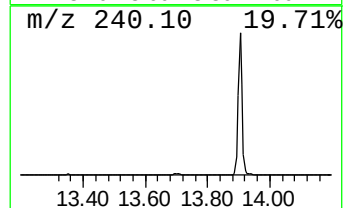
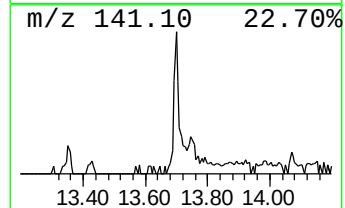
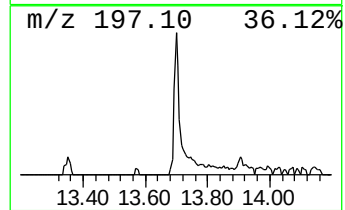
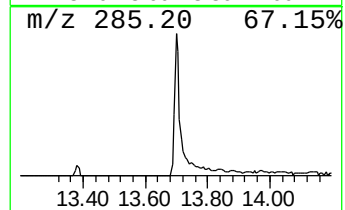
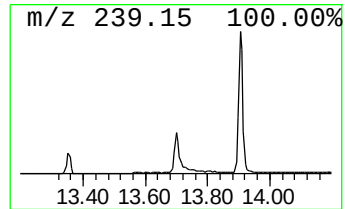
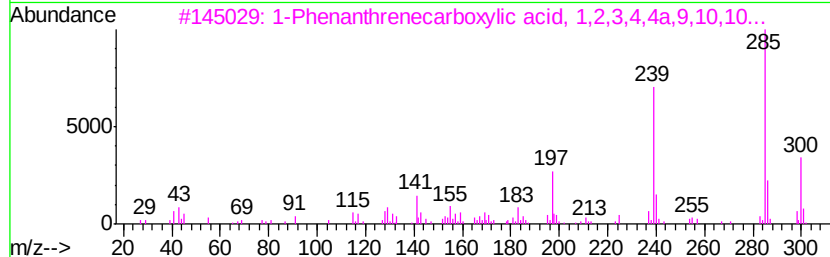
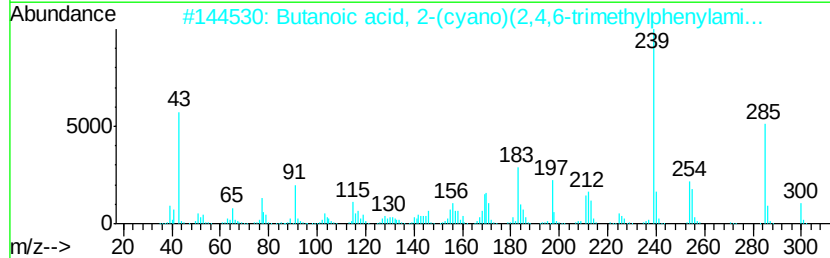
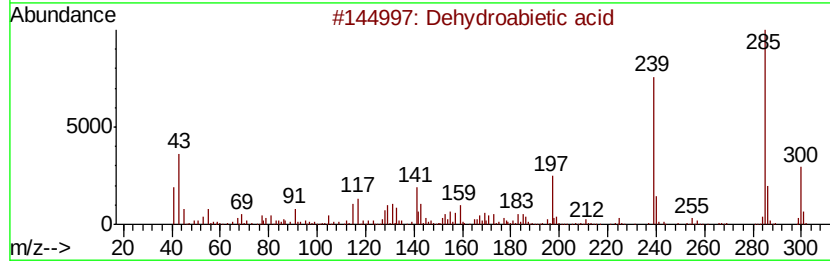
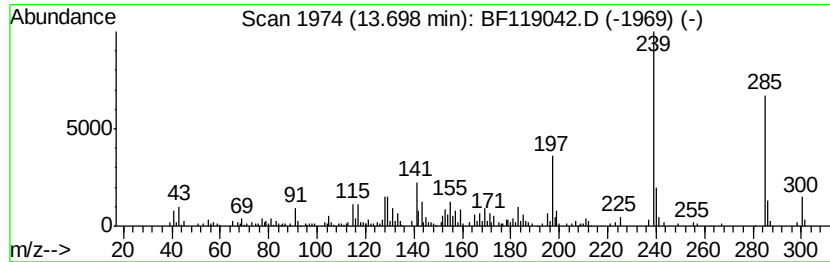
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Dehydroabiatic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.70	4.20 ng	240953	Chrysene-d12	13.90		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dehydroabiatic acid	300	C20H28O2	001740-19-8	95
2		Butanoic acid, 2-(cyano)(2,4,6-t...	300	C17H20N2O3	1000267-71-7	90
3		1-Phenanthrenecarboxylic acid, 1...	300	C20H28O2	005155-70-4	64
4		Ethyl 4-hydroxy-7-trifluoromethy...	285	C13H10F3N03	000391-02-6	58
5		3-Quinolinecarboxylic acid, 4-hy...	285	C13H10F3N03	023851-84-5	53



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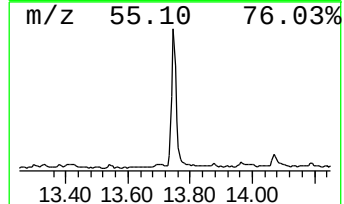
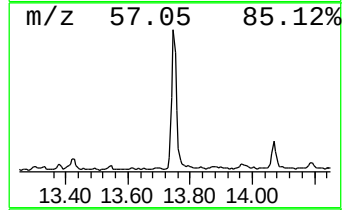
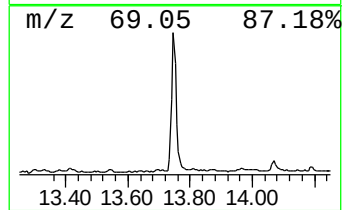
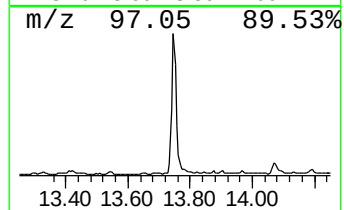
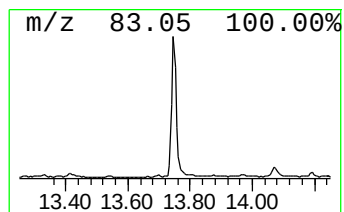
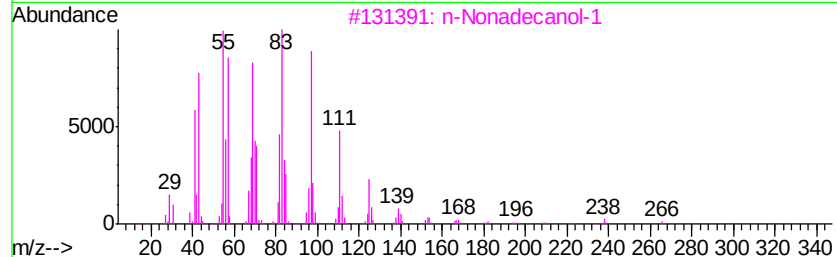
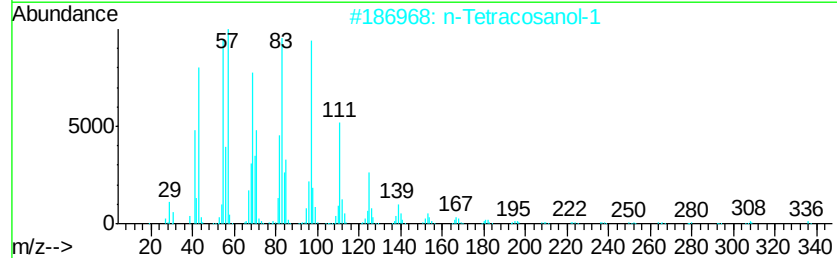
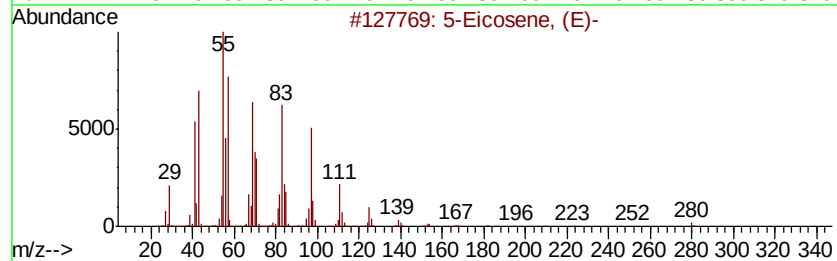
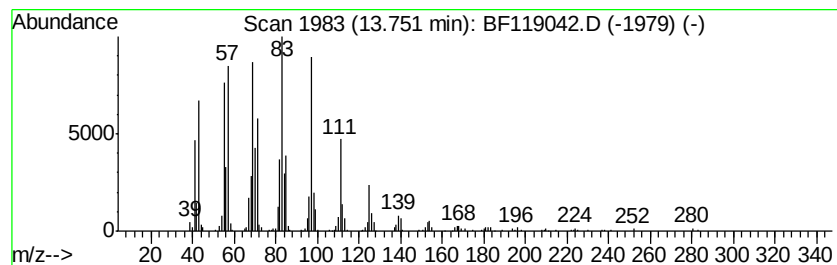
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 5-Eicosene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	12.59 ng	722459	Chrysene-d12	13.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2		n-Tetracosanol-1	354	C24H50O	000506-51-4	93
3		n-Nonadecanol-1	284	C19H40O	001454-84-8	93
4		Behenic alcohol	326	C22H46O	000661-19-8	91
5		1-Heneicosanol	312	C21H44O	015594-90-8	91



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Operator : CG/JU
Sample : L1635-10
Misc :
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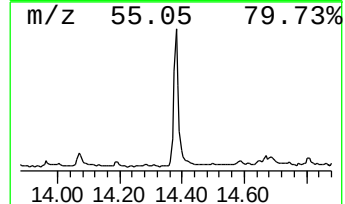
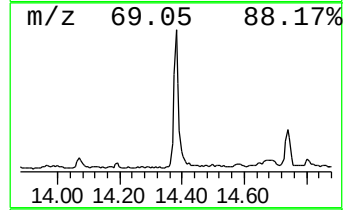
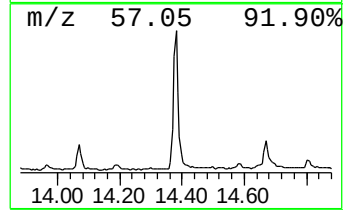
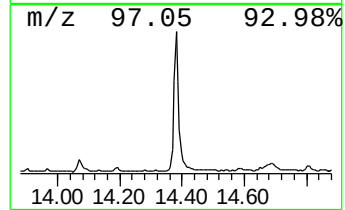
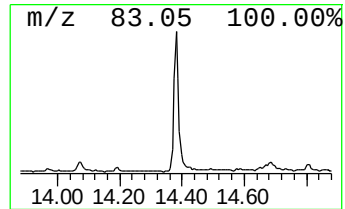
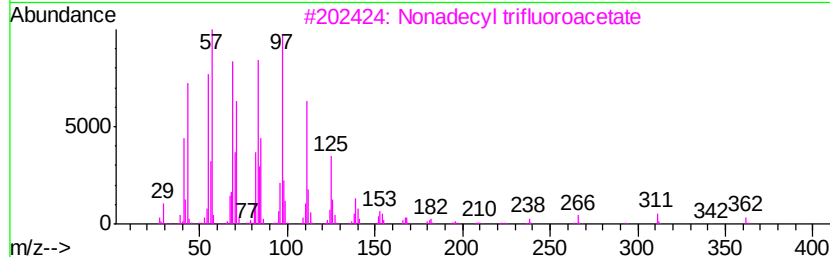
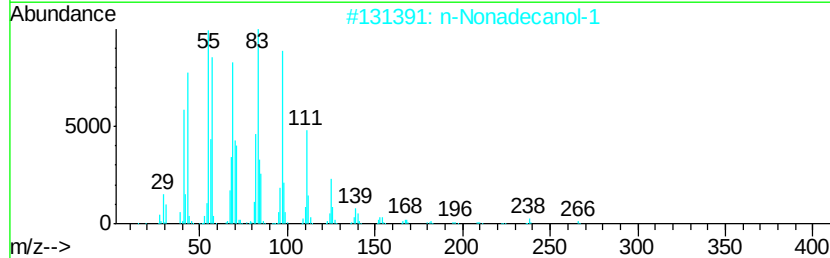
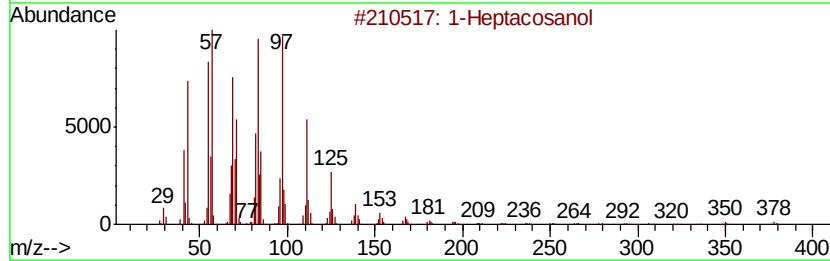
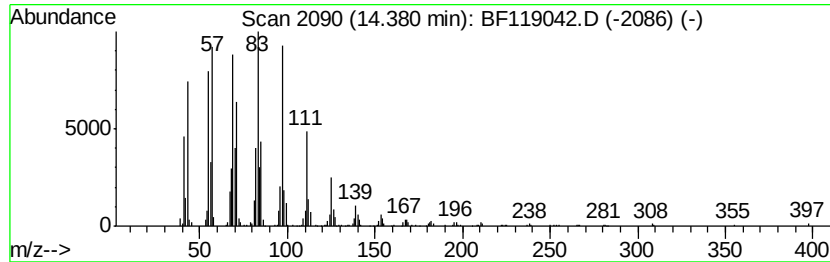
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Heptacosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.38	13.24 ng	759534	Chrysene-d12	13.90	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	95
2	n-Nonadecanol-1	284	C19H40O	001454-84-8	93
3	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	93
4	Behenic alcohol	326	C22H46O	000661-19-8	93
5	1-Octadecanol	270	C18H38O	000112-92-5	90



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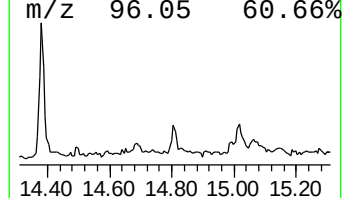
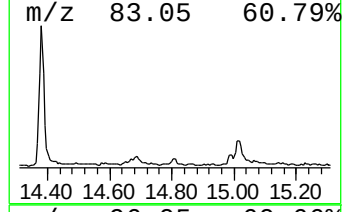
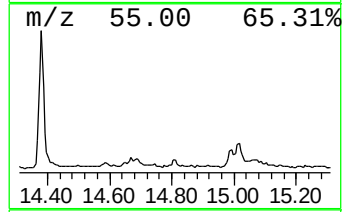
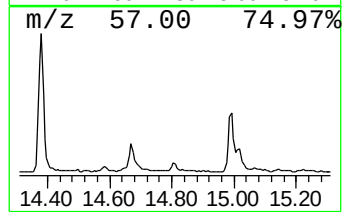
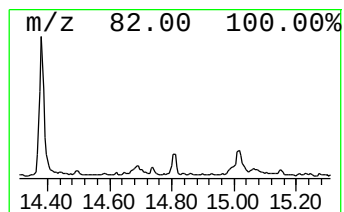
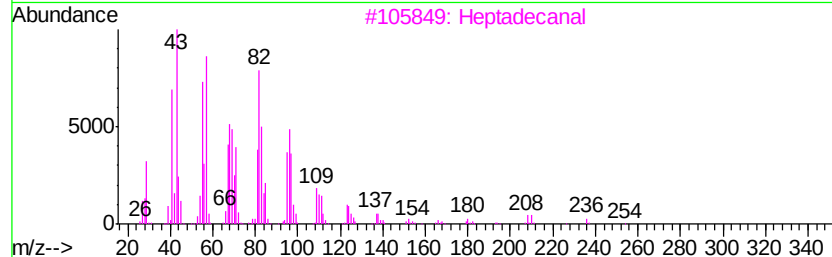
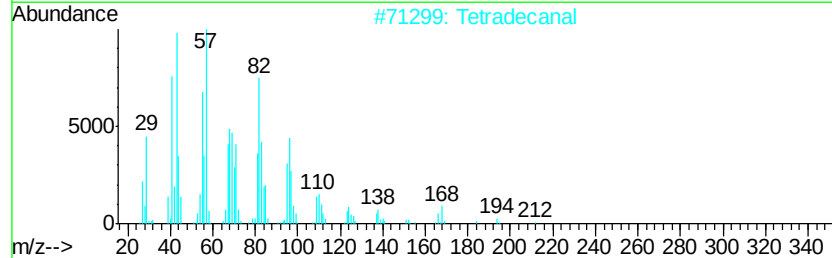
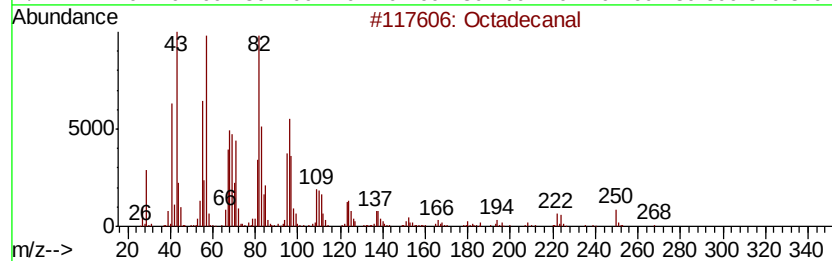
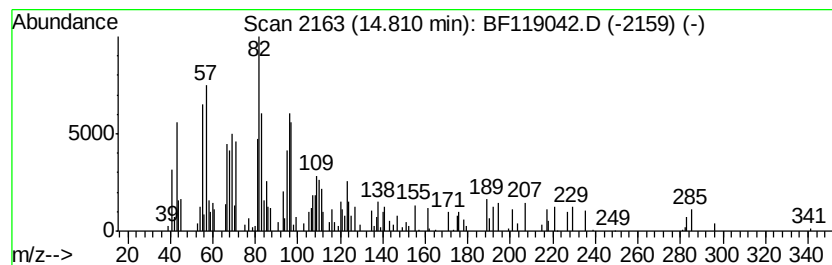
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanal Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	2.10 ng	116591	Perylene-d12	15.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanal	268	C18H36O	000638-66-4	83
2		Tetradecanal	212	C14H28O	000124-25-4	80
3		Heptadecanal	254	C17H34O	1000376-70-0	74
4		18-Nonadecen-1-ol	282	C19H38O	1000142-89-2	64
5		Cyclododecanol	184	C12H24O	001724-39-6	64



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Sample : L1635-10
Misc :
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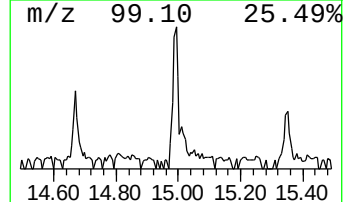
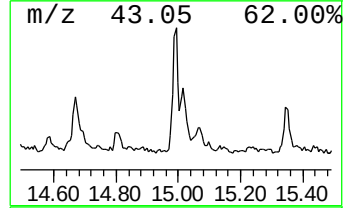
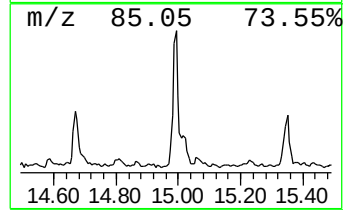
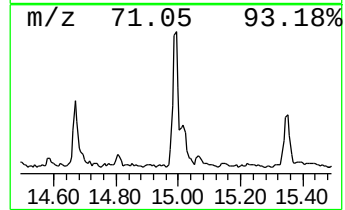
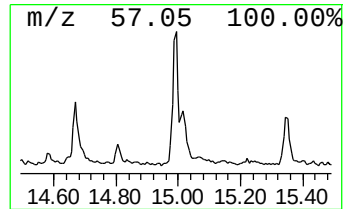
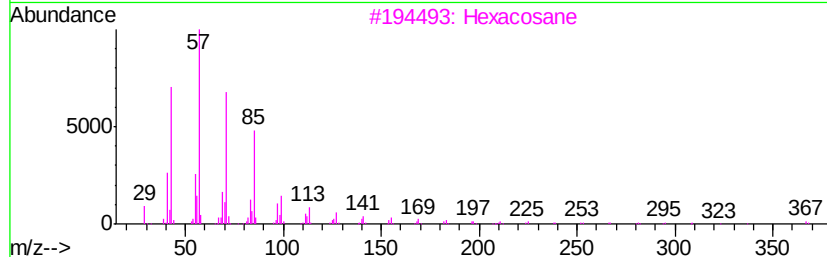
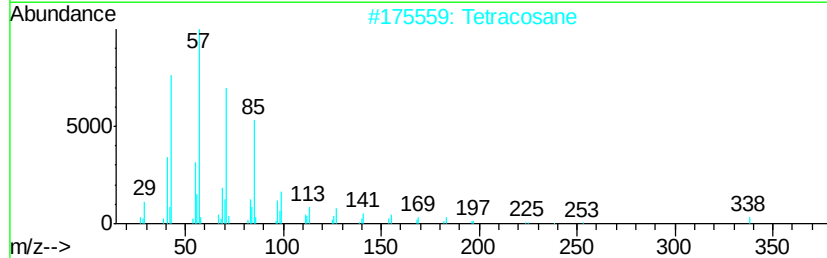
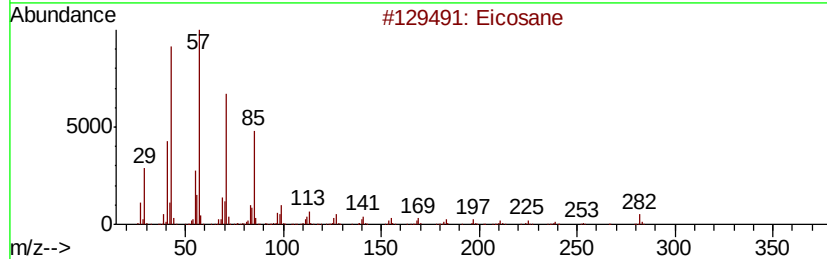
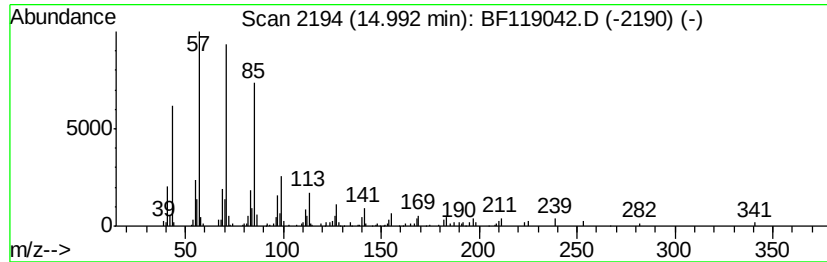
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.99	3.39 ng	188087	Perylene-d12	15.33
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane	282 C20H42	000112-95-8	93
2	Tetracosane	338 C24H50	000646-31-1	91
3	Hexacosane	366 C26H54	000630-01-3	91
4	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
5	Heneicosane, 11-(1-ethylpropyl)-	366 C26H54	055282-11-6	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022420\
Data File : BF119042.D
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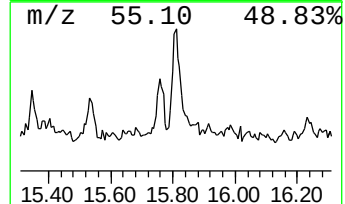
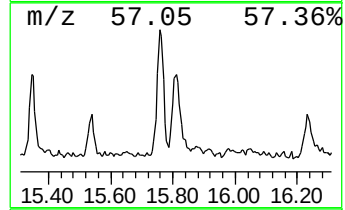
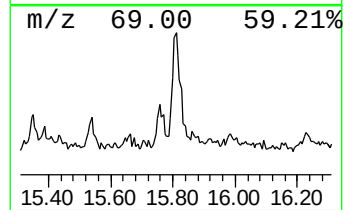
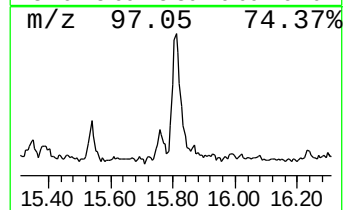
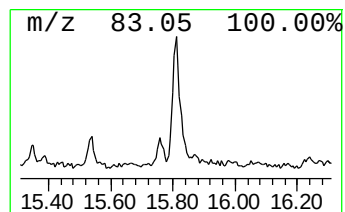
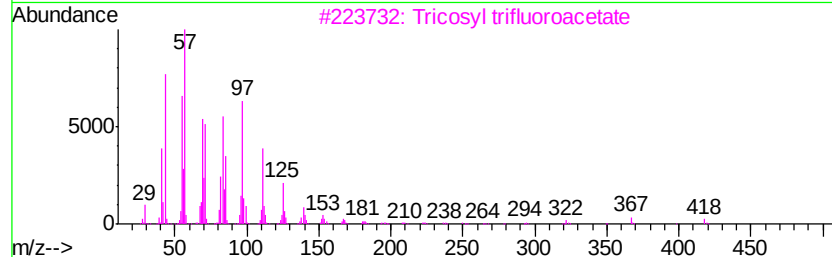
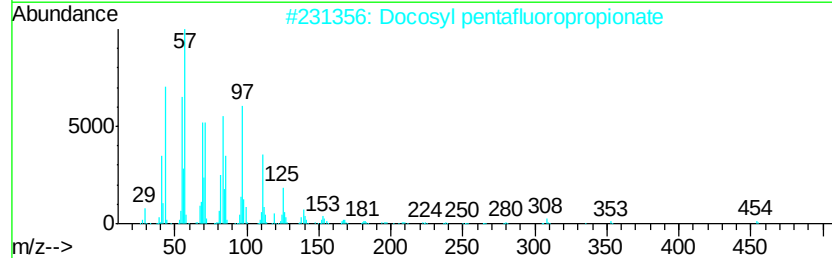
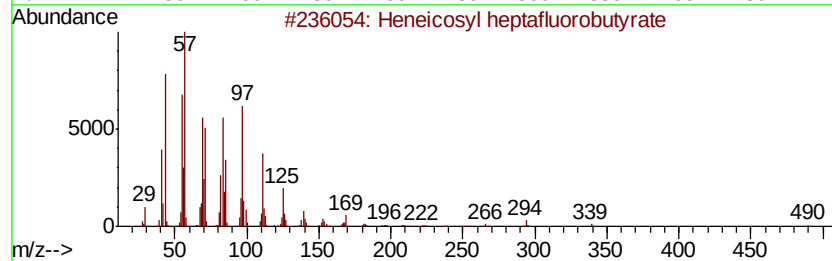
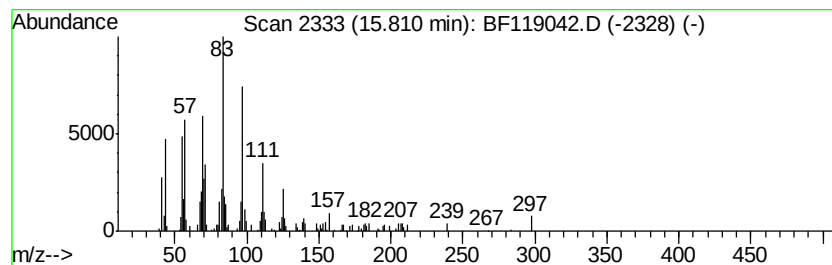
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Heneicosyl heptafluorobutyrate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.81	3.09 ng	170978	Perylene-d12	15.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosyl heptafluorobutyrate	508	C25H43F7O2	1000351-83-8	70
2		Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	62
3		Tricosyl trifluoroacetate	436	C25H47F3O2	1000351-75-1	62
4		Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	58
5		Octacosanol	410	C28H58O	000557-61-9	58



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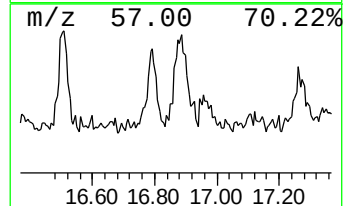
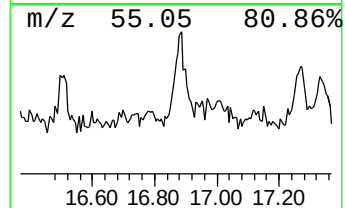
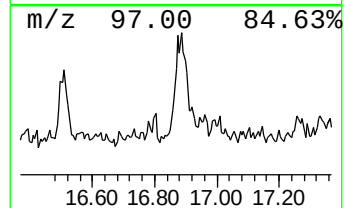
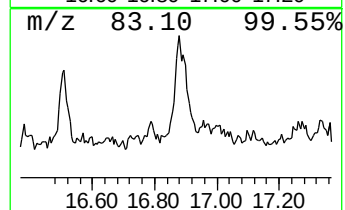
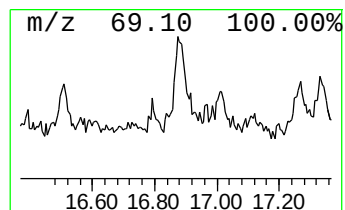
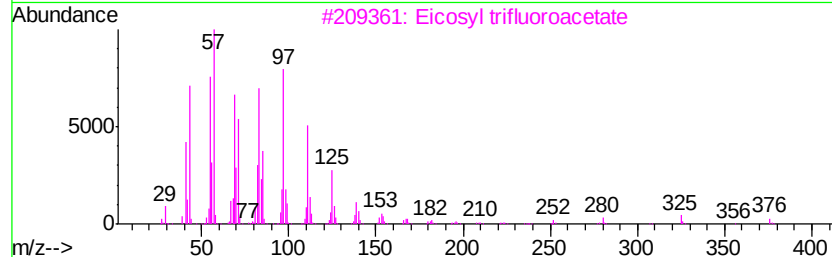
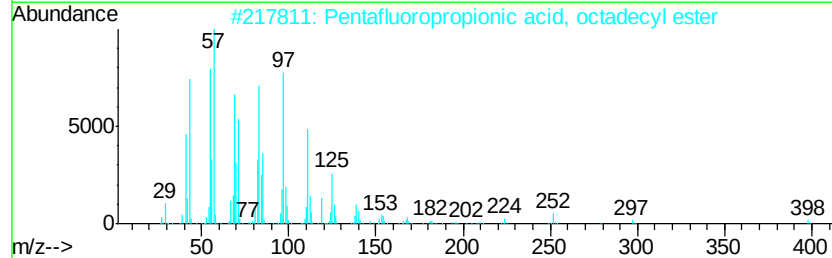
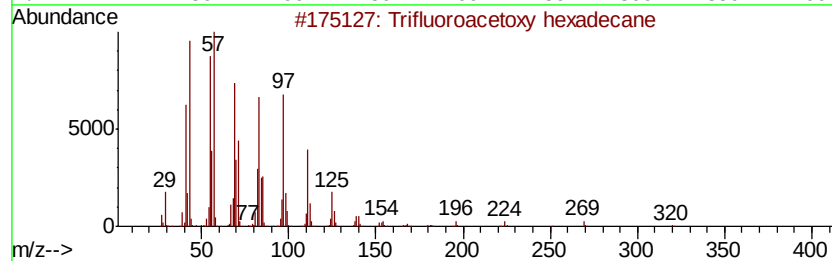
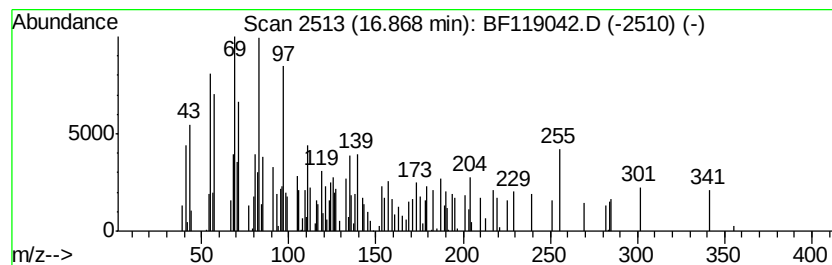
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Trifluoroacetoxy hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.87	2.41 ng	133649	Perylene-d12	15.33
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Trifluoroacetoxy hexadecane	338 C18H33F3O2	006222-03-3 72
2		Pentafluoropropionic acid, octad...	416 C21H37F5O2	959261-25-7 60
3		Eicosyl trifluoroacetate	394 C22H41F3O2	1000351-75-2 60
4		Heptadecyl trifluoroacetate	352 C19H35F3O2	1000351-87-0 53
5		1-Hexadecanesulfonyl chloride	324 C16H33ClO2S	038775-38-1 53



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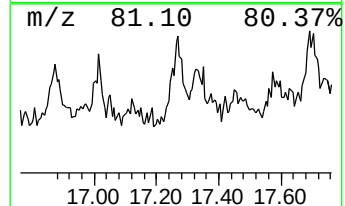
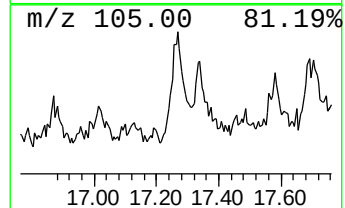
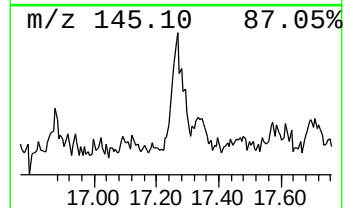
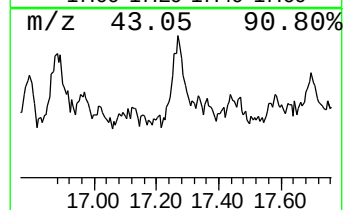
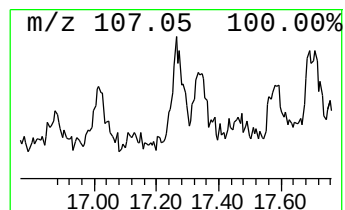
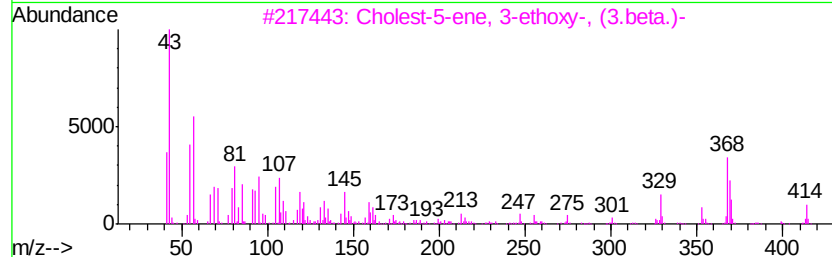
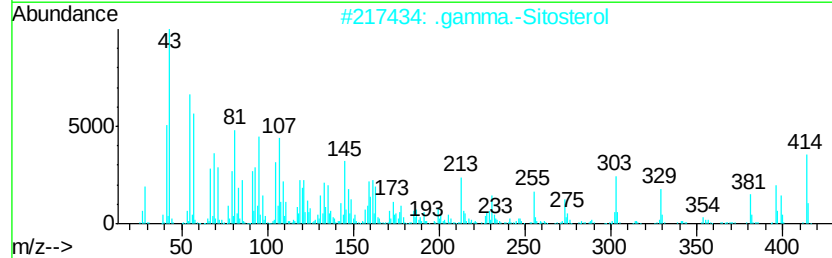
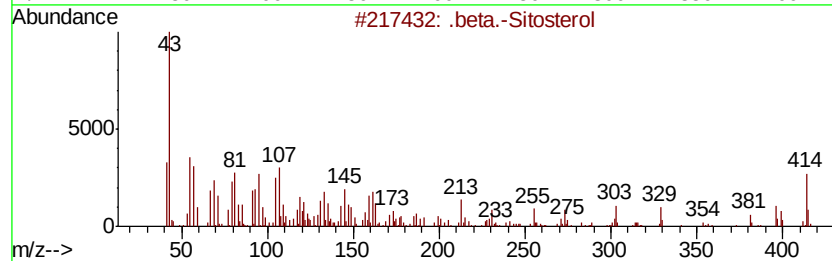
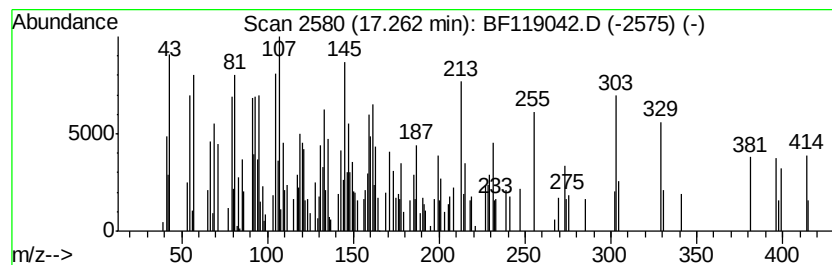
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 .beta.-Sitosterol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.26	3.31 ng	183204	Perylene-d12	15.33
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		.beta.-Sitosterol	414 C29H50O	000083-46-5 99
2		.gamma.-Sitosterol	414 C29H50O	000083-47-6 99
3		Cholest-5-ene, 3-ethoxy-, (3.beta...	414 C29H50O	000986-19-6 46
4		17-(1,5-Dimethylhexyl)-10,13-dim...	414 C29H50O	1000210-86-9 25
5		Stigmastan-3,5-diene	396 C29H48	1000214-16-4 22



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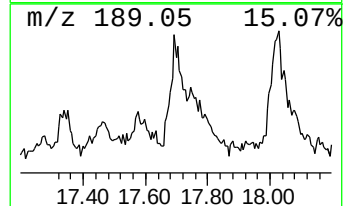
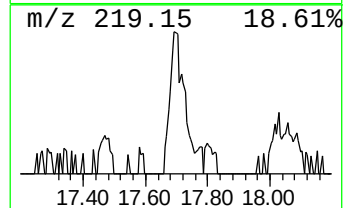
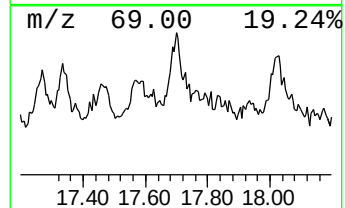
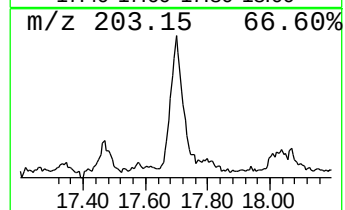
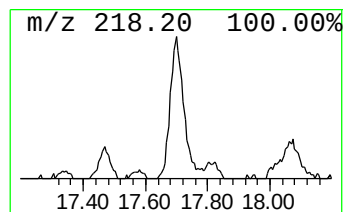
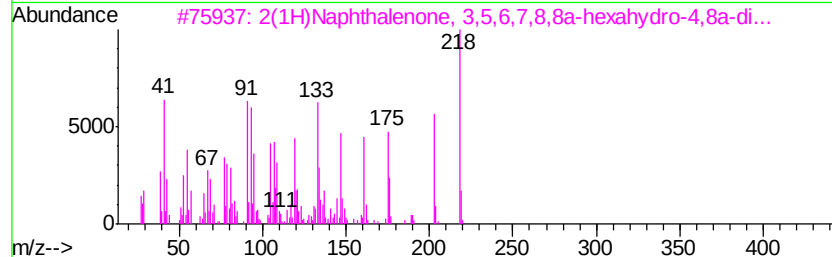
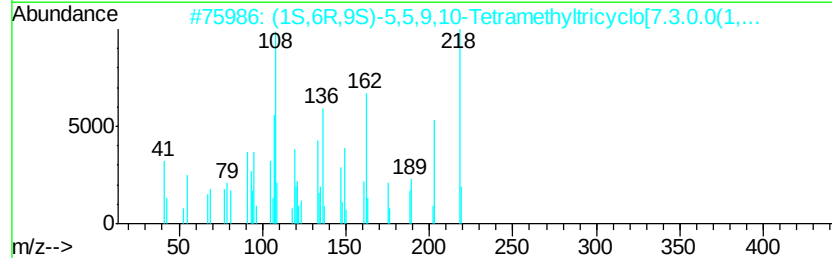
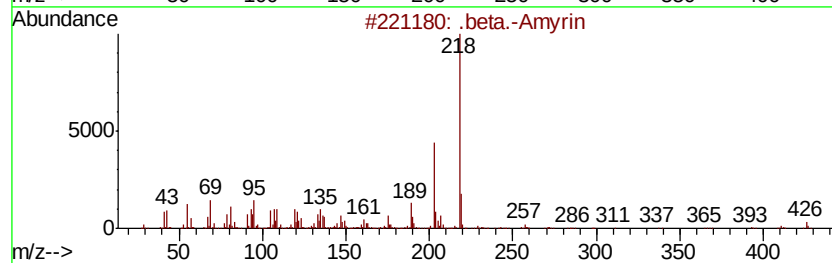
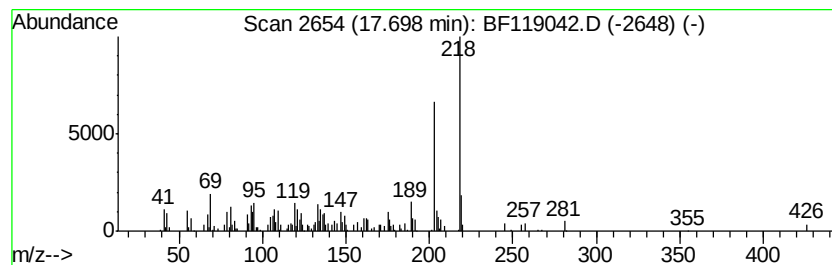
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 .beta.-Amyrin Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.70	3.51 ng	194245	Perylene-d12	15.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-Amyrin	426	C30H500	000559-70-6	95
2		(1S,6R,9S)-5,5,9,10-Tetramethylt...	218	C16H26	1000298-97-8	87
3		2(1H)Naphthalenone, 3,5,6,7,8,8a...	218	C15H220	1000188-66-5	81
4		Olean-12-ene	410	C30H50	000471-68-1	80
5		.beta.-Amyrin trimethylsilyl ether	498	C33H580Si	001721-67-1	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022420\
Data File : BF119042.D
Acq On : 25 Feb 2020 1:25
Operator : CG/JU
Sample : L1635-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

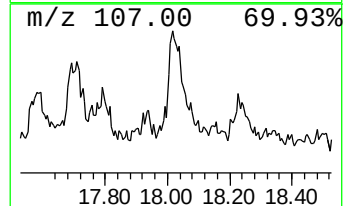
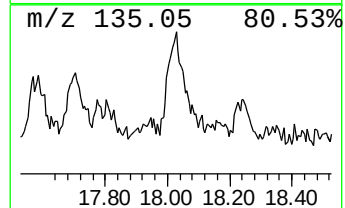
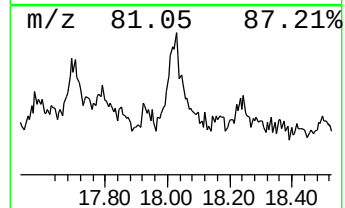
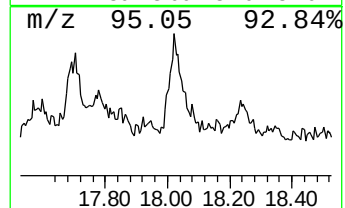
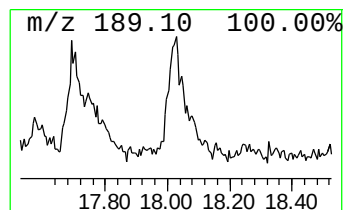
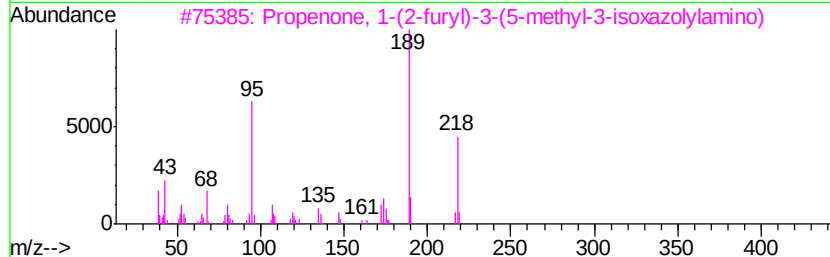
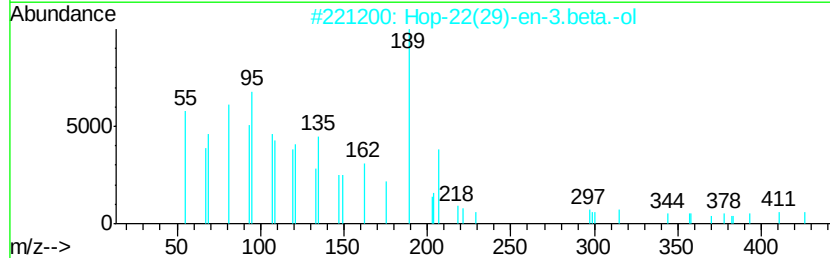
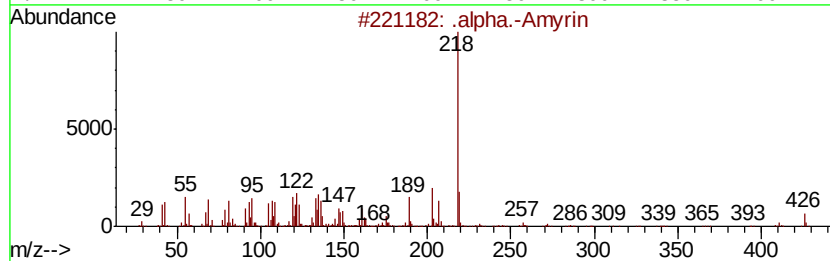
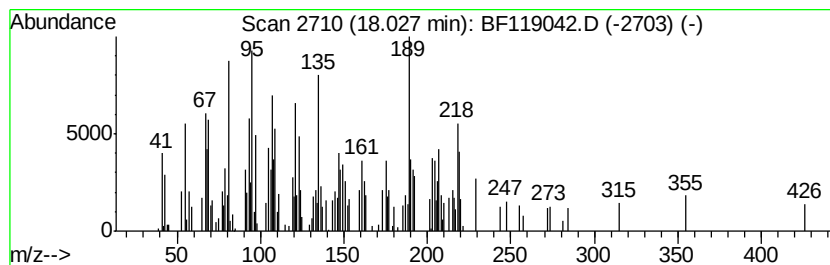
Instrument :
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ClientSampleId :
DUNR-TS007-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown18.03 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.03	4.55 ng	252399	Perylene-d12	15.33	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Amyrin	426	C30H50O	000638-95-9	46
2	Hop-22(29)-en-3.beta.-ol	426	C30H50O	058801-23-3	41
3	Propenone, 1-(2-furyl)-3-(5-meth...	218	C11H10N2O3	186957-33-5	35
4	Bicyclo[3.3.0]octan-3-one, 6-exo...	216	C9H13BrO	1000149-28-6	25
5	3-Amino-5-piperonyl-1,2,4-triazole	218	C10H10N4O2	014730-92-8	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022420\
Data File : BF119042.D
Acq On : 25 Feb 2020 1:25
Operator : CG/JU
Sample : L1635-10
Misc :
ALS Vial : 17 Sample Multiplier: 1

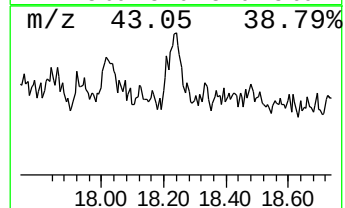
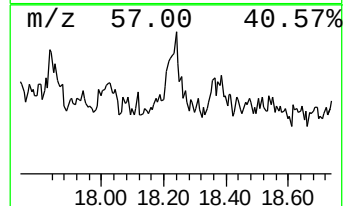
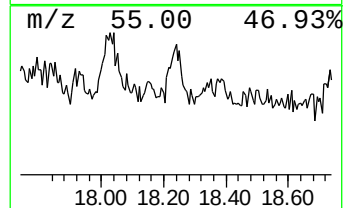
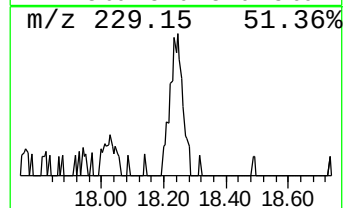
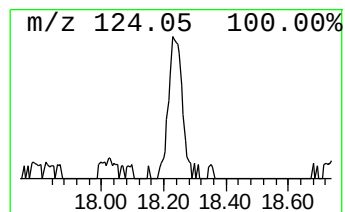
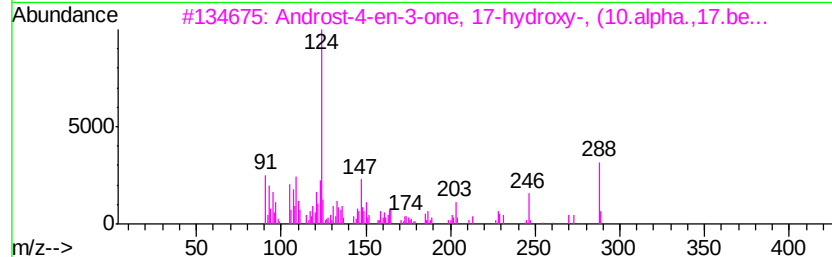
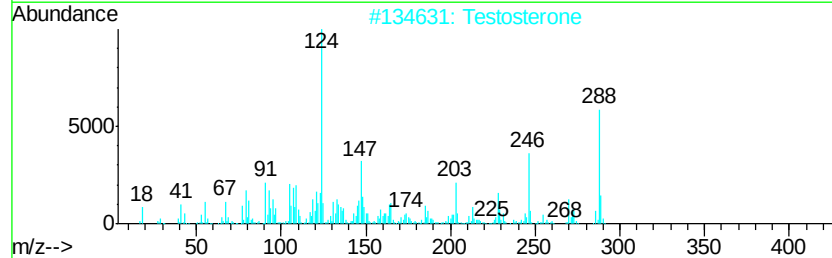
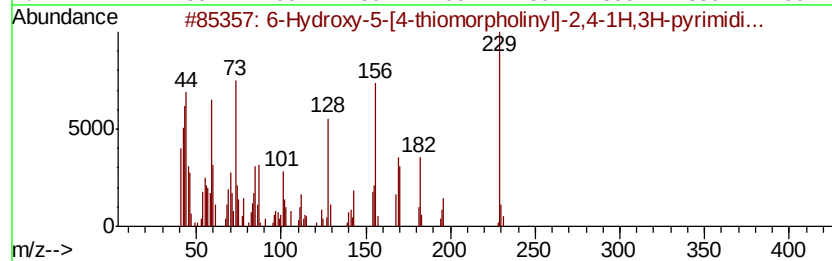
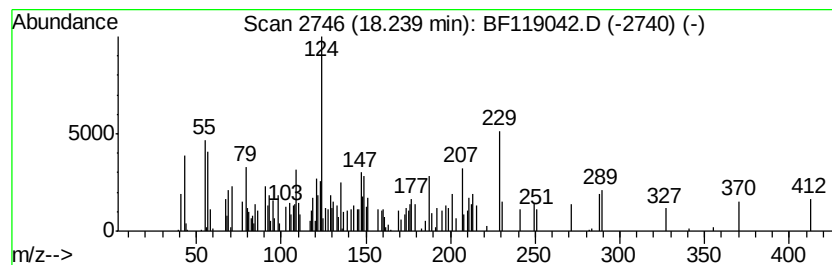
Instrument :
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ClientSampleId :
DUNR-TS007-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 6-Hydroxy-5-[4-thiomorpholi... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.24	2.08 ng	115415	Perylene-d12	15.33		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Hydroxy-5-[4-thiomorpholinyl]-...	229	C8H11N3O3S	1000212-11-4	53	
2	Testosterone	288	C19H28O2	000058-22-0	46	
3	Androst-4-en-3-one, 17-hydroxy-...	288	C19H28O2	000604-39-7	46	
4	3-Chloropropionic acid, 2,6-dime...	256	C14H21ClO2	1000299-21-8	25	
5	Benzeneacetic acid, .alpha.-meth...	271	C17H21NO2	000500-55-0	25	



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 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUNR-TS007-01

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.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
n-Hexadecanoic acid	11.80	6.4 ng		517069	4	11.27	1627120	20.0
10,18-Bisnorabiet...	12.52	5.1 ng		416347	4	11.27	1627120	20.0
Octadecanoic acid	12.58	2.1 ng		172005	4	11.27	1627120	20.0
Dehydroabietic acid	13.70	4.2 ng		240953	5	13.90	1147710	20.0
5-Eicosene, (E)-	13.75	12.6 ng		722459	5	13.90	1147710	20.0
1-Heptacosanol	14.38	13.2 ng		759534	5	13.90	1147710	20.0
Octadecanal	14.81	2.1 ng		116591	6	15.33	1108280	20.0
Eicosane	14.99	3.4 ng		188087	6	15.33	1108280	20.0
Heneicosyl heptaf...	15.81	3.1 ng		170978	6	15.33	1108280	20.0
Trifluoroacetoxy ...	16.87	2.4 ng		133649	6	15.33	1108280	20.0
.beta.-Sitosterol	17.26	3.3 ng		183204	6	15.33	1108280	20.0
.beta.-Amyrin	17.70	3.5 ng		194245	6	15.33	1108280	20.0
unknown18.03	18.03	4.5 ng		252399	6	15.33	1108280	20.0
6-Hydroxy-5-[4-th...	18.24	2.1 ng		115415	6	15.33	1108280	20.0