

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
 Data File : BF121086.D
 Acq On : 28 Jul 2020 23:58
 Operator : JU/CG
 Sample : L3435-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-2

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-BF071620.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	5.040	499	502	509	rVB	40281	56530	1.03%	0.136%
2	5.434	563	569	572	rBV	3625091	3966223	72.40%	9.575%
3	6.451	736	742	745	rBV	3348470	3941022	71.94%	9.514%
4	6.645	772	775	782	rBV	49846	57730	1.05%	0.139%
5	6.810	800	803	807	rBV	1130298	893614	16.31%	2.157%
6	6.875	811	814	827	rBV	56126	87225	1.59%	0.211%
7	7.381	894	900	903	rBV	2497398	2721877	49.68%	6.571%
8	8.092	1017	1021	1025	rBV	1284995	1148901	20.97%	2.773%
9	9.175	1200	1205	1208	rBV	4191077	4556983	83.18%	11.001%
10	9.486	1254	1258	1259	rBV	83832	74430	1.36%	0.180%
11	9.504	1259	1261	1265	rVB	116125	93574	1.71%	0.226%
12	9.563	1265	1271	1280	rBV	625402	560437	10.23%	1.353%
13	9.845	1316	1319	1323	rBV	1414551	1307722	23.87%	3.157%
14	9.875	1323	1324	1337	rVB2	77262	85596	1.56%	0.207%
15	10.057	1350	1355	1373	rBV2	289282	367982	6.72%	0.888%
16	10.522	1425	1434	1437	rBV	242456	223999	4.09%	0.541%
17	10.639	1447	1454	1468	rBV	2577353	2978651	54.37%	7.191%
18	10.745	1470	1472	1478	rVB	54039	79407	1.45%	0.192%
19	10.998	1510	1515	1526	rBV2	102721	141636	2.59%	0.342%
20	11.333	1568	1572	1574	rBV	1669371	1541359	28.14%	3.721%
21	11.351	1574	1575	1579	rVB	153795	120059	2.19%	0.290%
22	11.863	1658	1662	1669	rBV2	363777	335444	6.12%	0.810%
23	11.945	1673	1676	1685	rVB3	47147	67499	1.23%	0.163%
24	12.310	1734	1738	1740	rBV	85842	80063	1.46%	0.193%
25	12.345	1740	1744	1748	rVV3	54541	61693	1.13%	0.149%
26	12.539	1773	1777	1780	rBV2	257045	244817	4.47%	0.591%
27	12.598	1783	1787	1792	rVV2	98343	131153	2.39%	0.317%
28	12.645	1792	1795	1801	rVV	108776	115159	2.10%	0.278%
29	12.768	1809	1816	1819	rBV	184582	197837	3.61%	0.478%
30	12.857	1828	1831	1834	rVB	113902	104356	1.90%	0.252%
31	12.921	1837	1842	1846	rBV	4288824	5478418	100.00%	13.225%
32	13.127	1874	1877	1879	rBV	73604	75532	1.38%	0.182%
33	13.345	1911	1914	1918	rBV	224779	219016	4.00%	0.529%
34	13.486	1934	1938	1941	rBV3	134867	135998	2.48%	0.328%

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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	13.815	1990	1994	2004	rBV	1360841	1681873	30.70%	4.060%
36	13.968	2015	2020	2023	rBV	1784411	1734960	31.67%	4.188%
37	14.133	2045	2048	2060	rVB2	102875	134093	2.45%	0.324%
38	14.404	2089	2094	2097	rBV	203344	288011	5.26%	0.695%
39	14.445	2097	2101	2108	rVB2	334704	486086	8.87%	1.173%
40	14.739	2147	2151	2154	rBV	84707	103696	1.89%	0.250%
41	14.886	2173	2176	2183	rVB	66118	89942	1.64%	0.217%
42	15.074	2204	2208	2210	rBV	282857	307606	5.61%	0.743%
43	15.104	2210	2213	2219	rVB	196191	269833	4.93%	0.651%
44	15.415	2261	2266	2276	rBV	1167859	1766705	32.25%	4.265%
45	15.639	2301	2304	2308	rVB	76398	95109	1.74%	0.230%
46	15.874	2339	2344	2348	rBV	317758	485362	8.86%	1.172%
47	15.927	2348	2353	2361	rBV	345076	666316	12.16%	1.609%
48	16.362	2423	2427	2431	rBV2	46413	79968	1.46%	0.193%
49	16.645	2471	2475	2481	rVB3	72785	129198	2.36%	0.312%
50	16.939	2521	2525	2532	rVB2	84168	154635	2.82%	0.373%
51	17.039	2535	2542	2547	rBV3	123105	364468	6.65%	0.880%
52	17.415	2600	2606	2616	rBV3	124343	334432	6.10%	0.807%

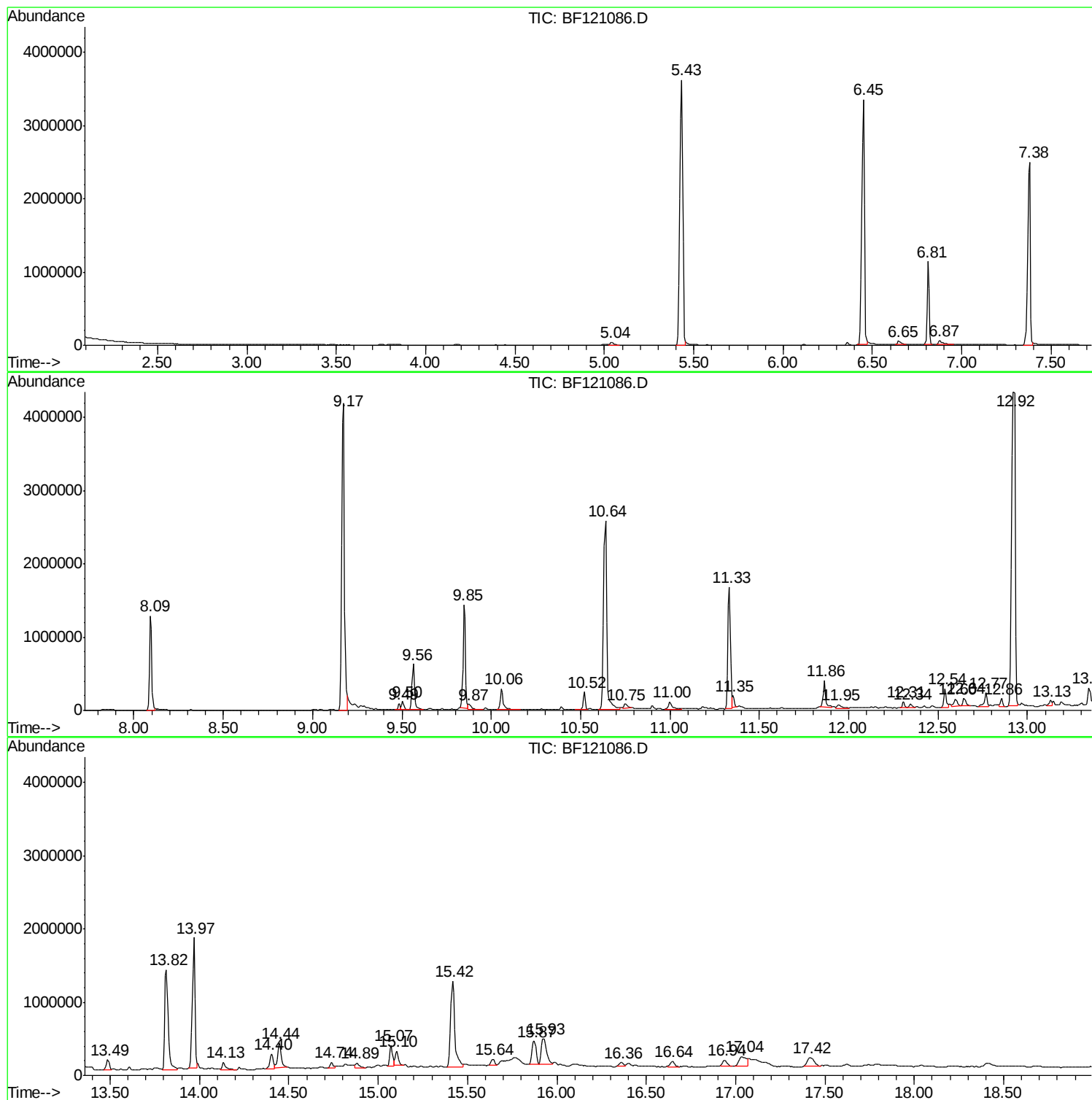
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ClientSampled :
SP-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.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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Instrument :
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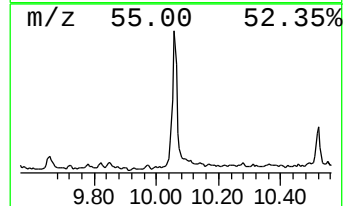
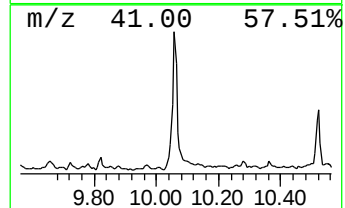
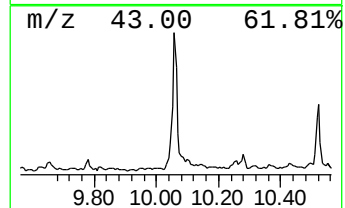
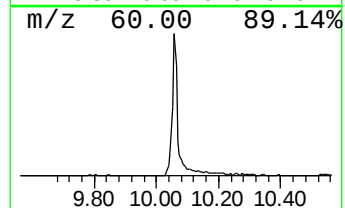
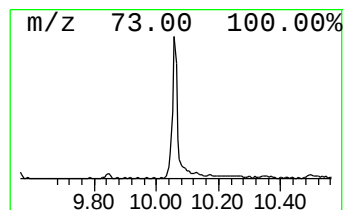
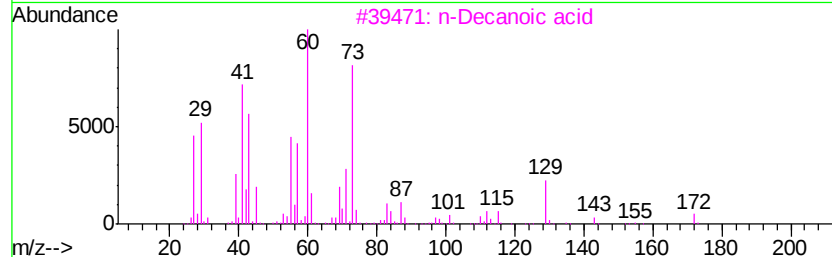
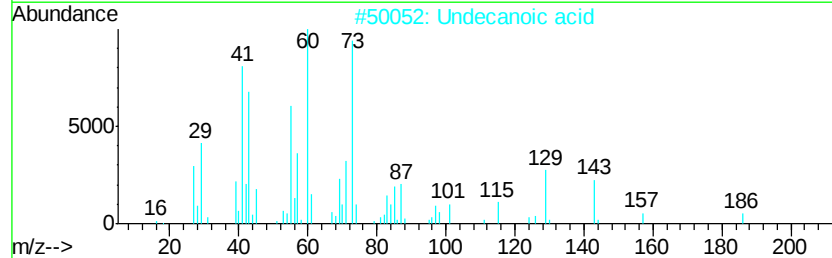
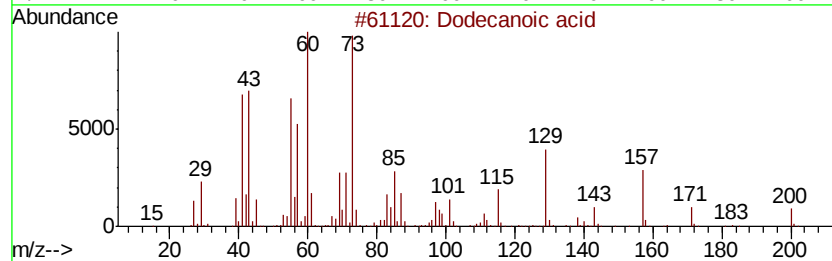
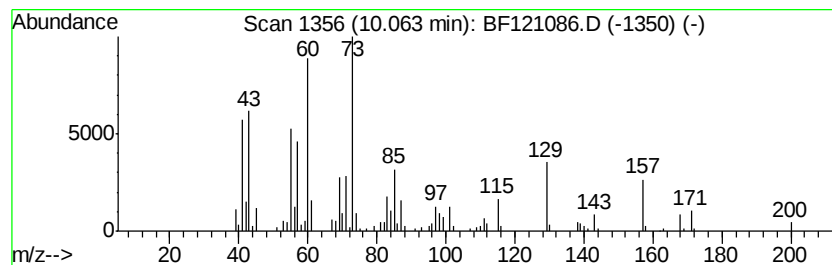
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Dodecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.06	5.63 ng	367982	Acenaphthene-d10	9.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecanoic acid	200	C12H24O2	000143-07-7	98
2		Undecanoic acid	186	C11H22O2	000112-37-8	72
3		n-Decanoic acid	172	C10H20O2	000334-48-5	53
4		Maltose	342	C12H22O11	000069-79-4	47
5		Octanoic acid	144	C8H16O2	000124-07-2	46



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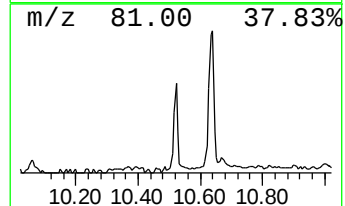
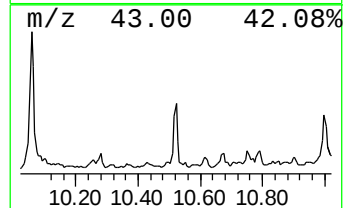
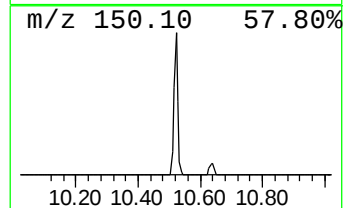
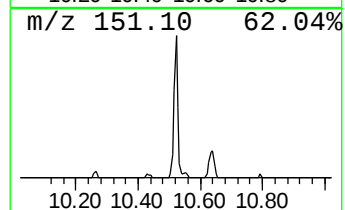
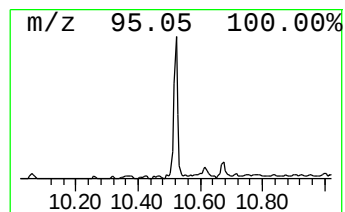
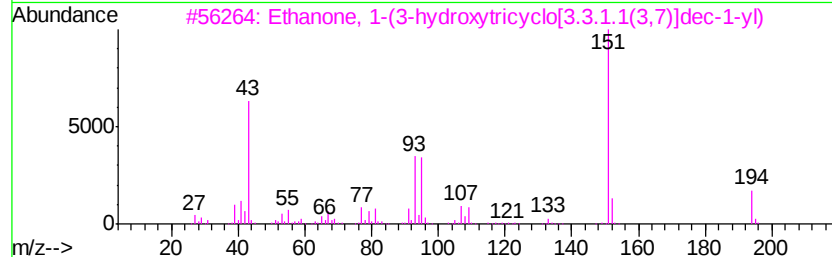
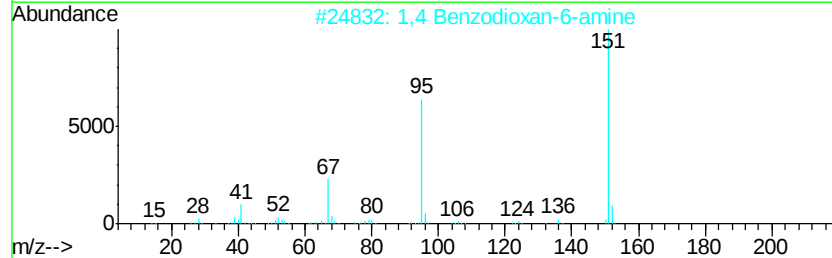
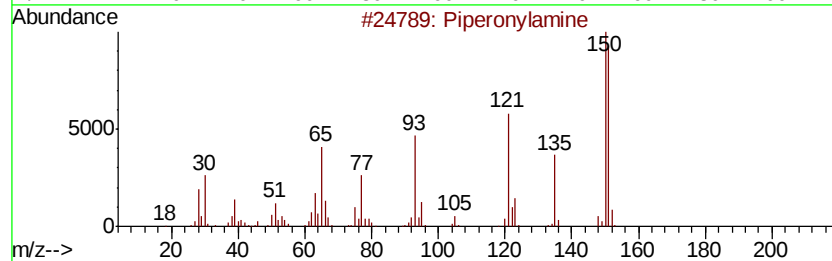
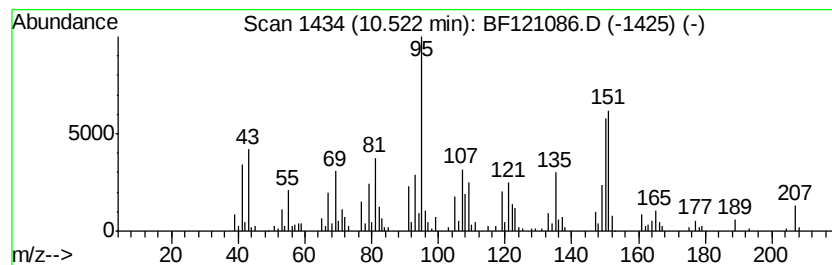
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown10.52 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.52	3.43 ng	223999	Acenaphthene-d10	9.85	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Piperonylamine	151	C8H9NO2	002620-50-0	38
2	1,4 Benzodioxan-6-amine	151	C8H9NO2	022013-33-8	30
3	Ethanone, 1-(3-hydroxytricyclo[3...]	194	C12H18O2	039917-38-9	27
4	2-Cyclopenten-1-one, 2-(2-butenyl...]	150	C10H14O	017190-71-5	25
5	Pyrazolo[1,5-a]pyridine, 3,3a,4,...]	150	C9H14N2	1000221-84-2	22



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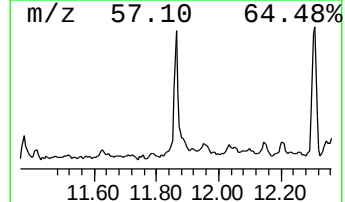
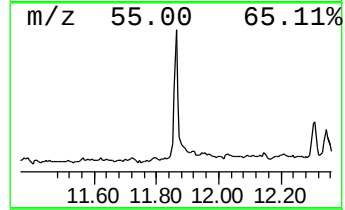
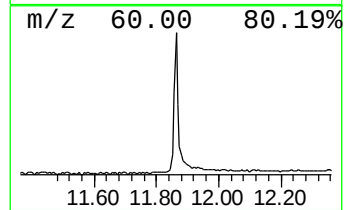
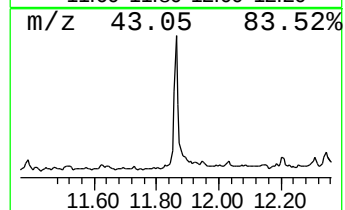
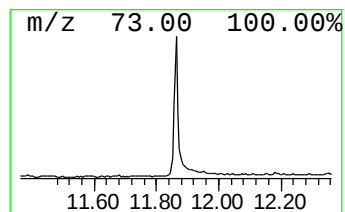
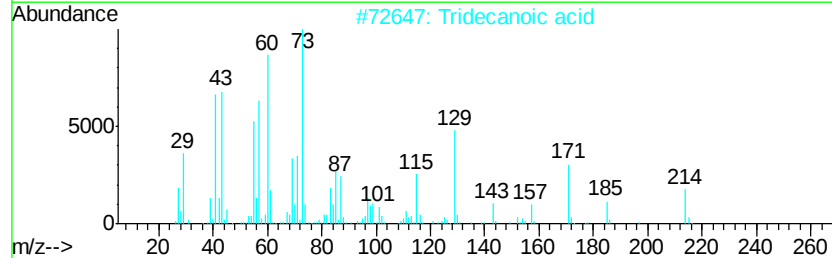
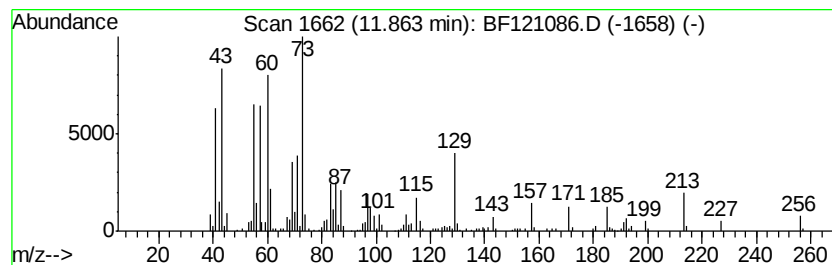
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	4.35 ng	335444	Phenanthrene-d10	11.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	94
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	18



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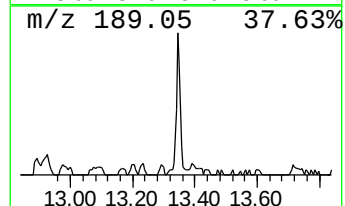
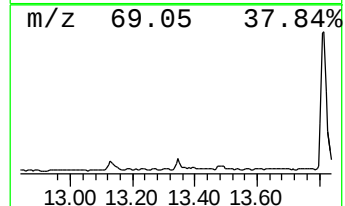
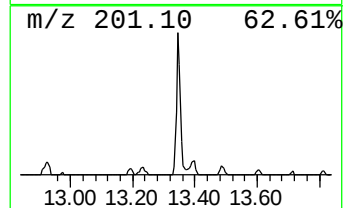
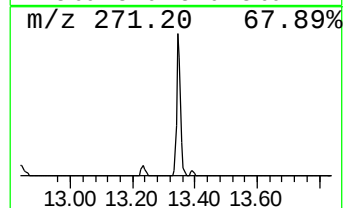
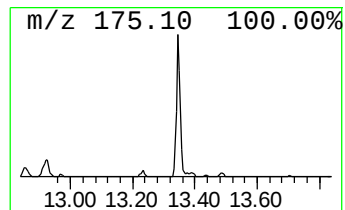
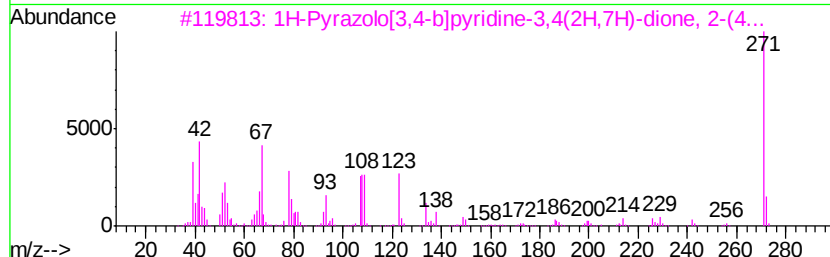
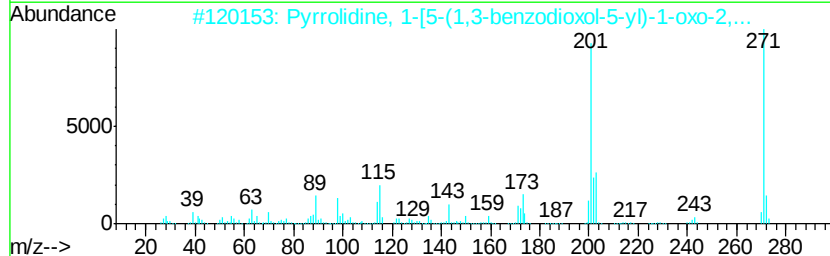
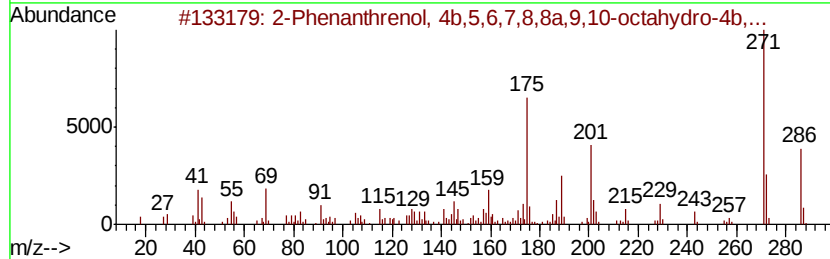
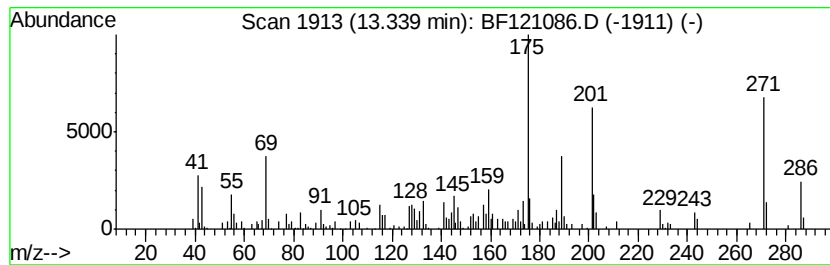
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.34	2.52 ng	219016	Chrysene-d12	13.97		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol,	4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	96
2	Pyrrolidine, 1-[5-(1,3-benzodiox...		271	C16H17N03	025924-78-1	46
3	1H-Pyrazolo[3,4-b]pyridine-3,4(2...		271	C13H13N5O2	1000362-36-1	18
4	Dihydronormorphinone		271	C16H17N03	014696-23-2	14
5	3-Cyano-6-trifluoromethylphenant...		271	C16H8F3N	1000253-90-3	14



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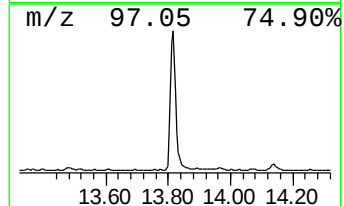
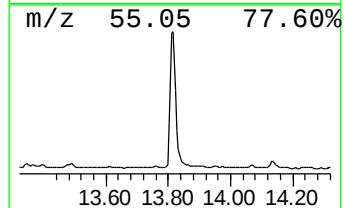
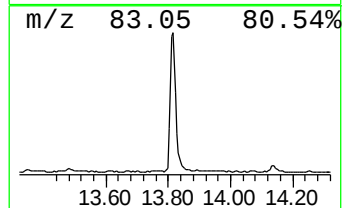
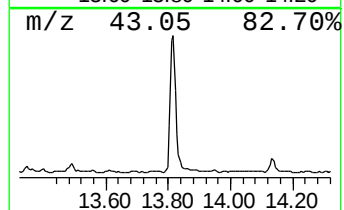
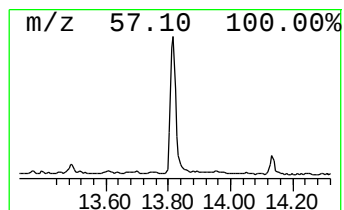
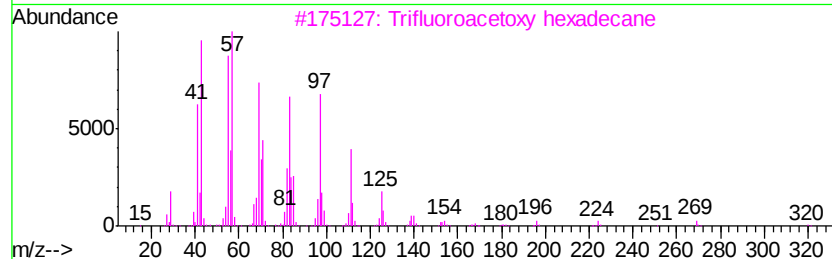
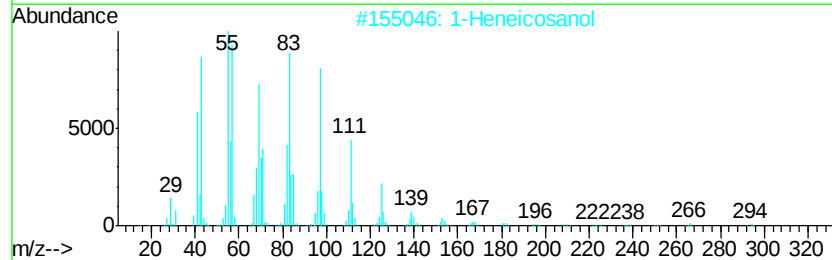
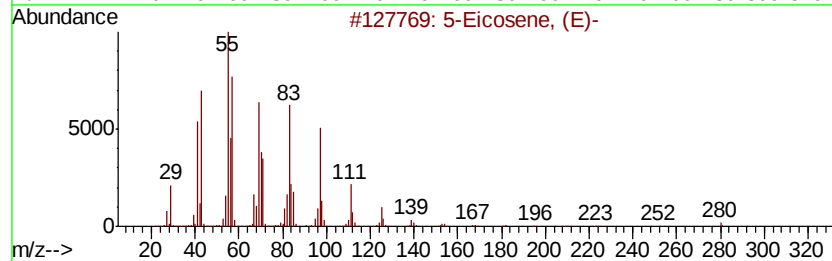
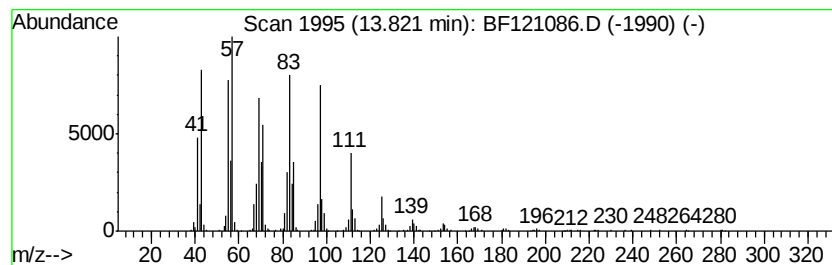
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 5-Eicosene, (E)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	19.39 ng	1681870	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2		1-Heneicosanol	312	C21H44O	015594-90-8	95
3		Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
4		Heptadecyl heptafluorobutyrat	452	C21H35F7O2	959085-66-6	94
5		Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121086.D
Acq On : 28 Jul 2020 23:58
Operator : JU/CG
Sample : L3435-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

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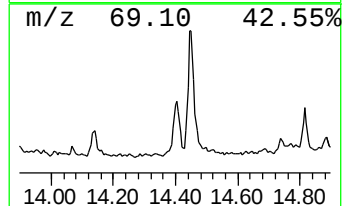
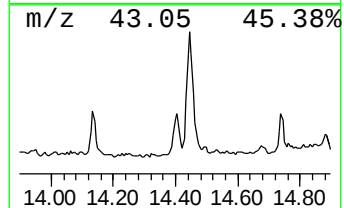
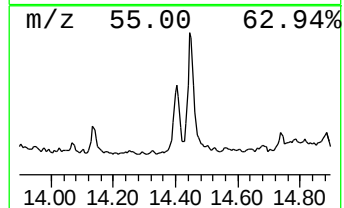
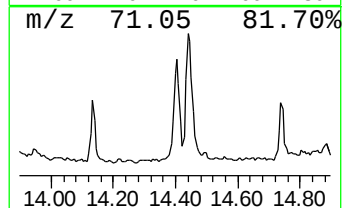
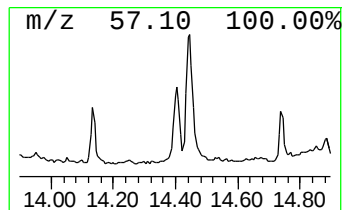
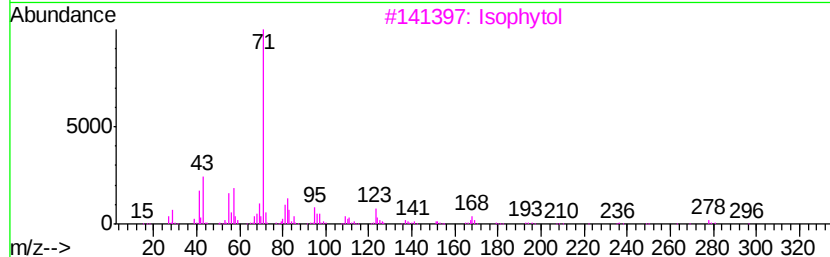
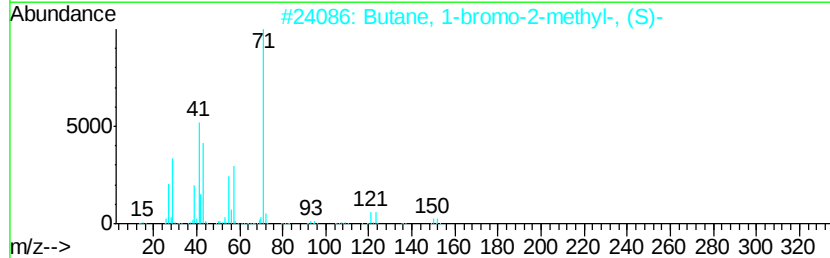
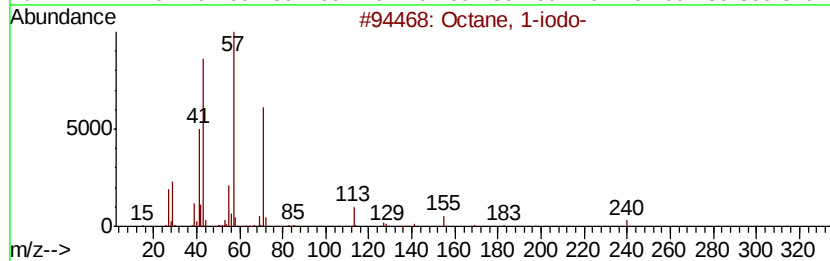
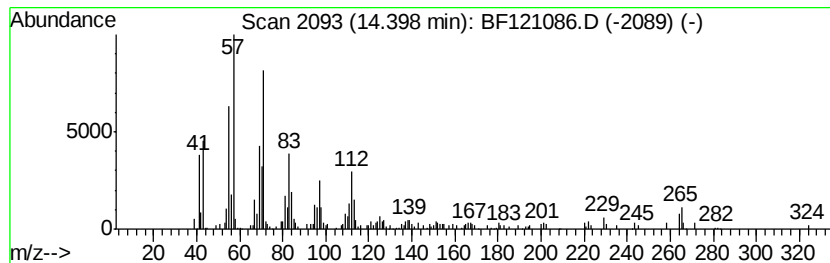
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown14.40 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.40	3.32 ng	288011	Chrysene-d12	13.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 1-iodo-	240	C8H17I	000629-27-6	38
2		Butane, 1-bromo-2-methyl-, (S)-	150	C5H11Br	000534-00-9	38
3		Isophytol	296	C20H40O	000505-32-8	35
4		3-Cyclopentylpropionic acid, 2-t...	226	C13H22O3	1000293-46-9	35
5		Z-5-Nonadecene	266	C19H38	1000131-11-8	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
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Sample : L3435-03
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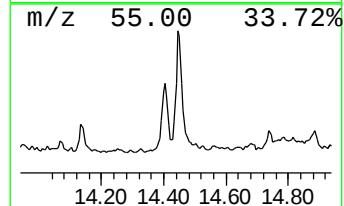
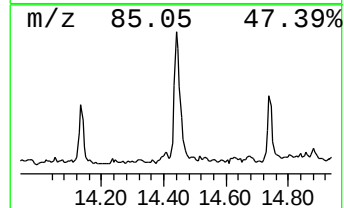
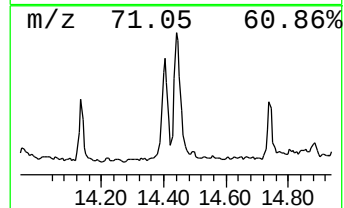
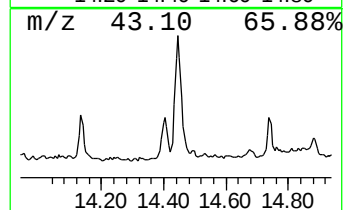
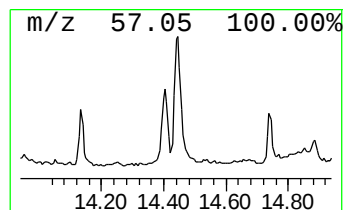
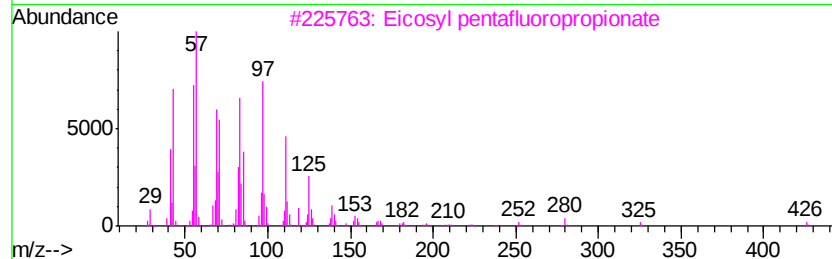
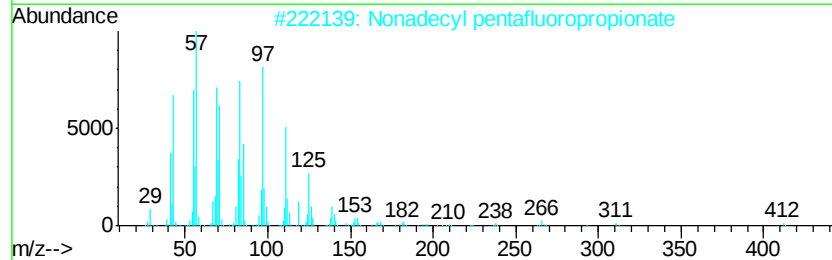
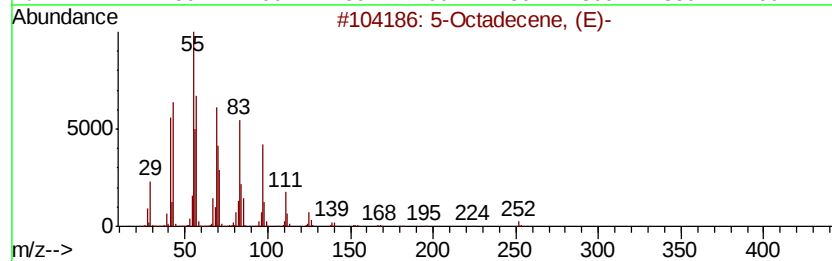
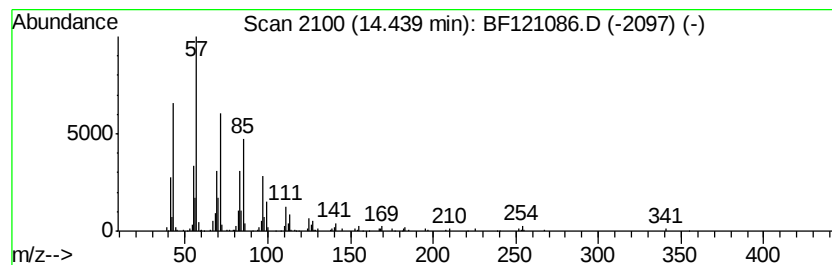
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 5-Octadecene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.44	5.60 ng	486086	Chrysene-d12	13.97		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Octadecene, (E)-		252	C18H36	007206-21-5	93
2	Nonadecyl pentafluoropropionate		430	C22H39F5O2	1000351-88-8	91
3	Eicosyl pentafluoropropionate		444	C23H41F5O2	1000351-80-8	91
4	Docosyl trifluoroacetate		422	C24H45F3O2	1000351-75-5	91
5	17-Pentatriacontene		491	C35H70	006971-40-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121086.D
Acq On : 28 Jul 2020 23:58
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Sample : L3435-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

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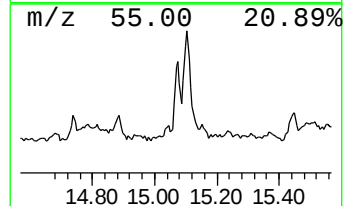
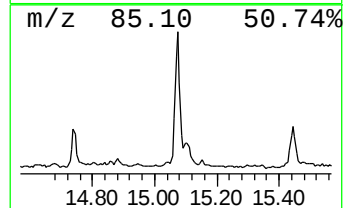
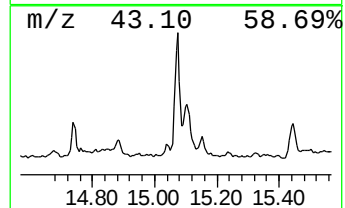
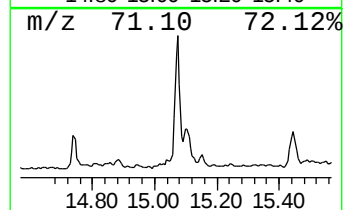
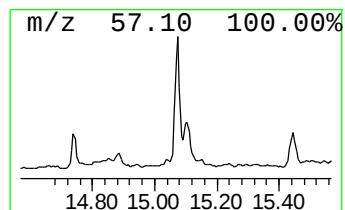
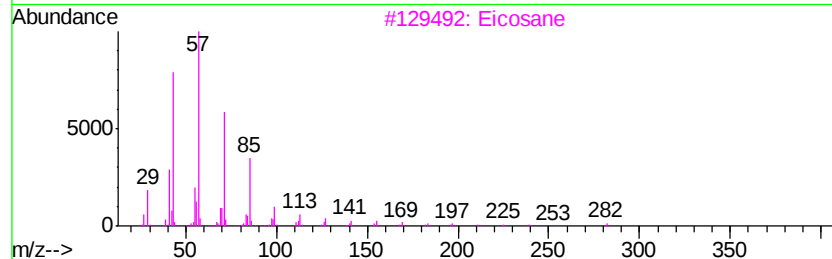
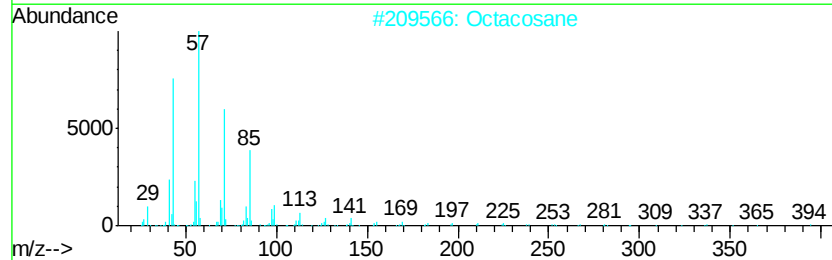
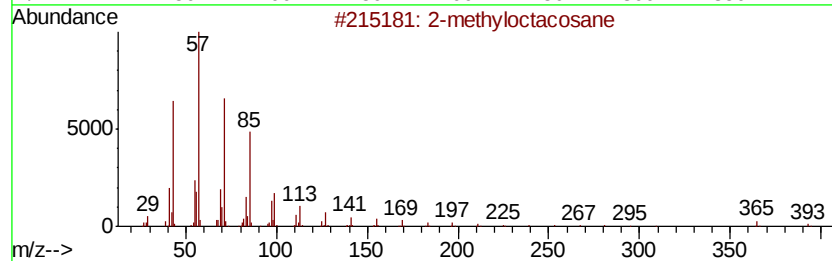
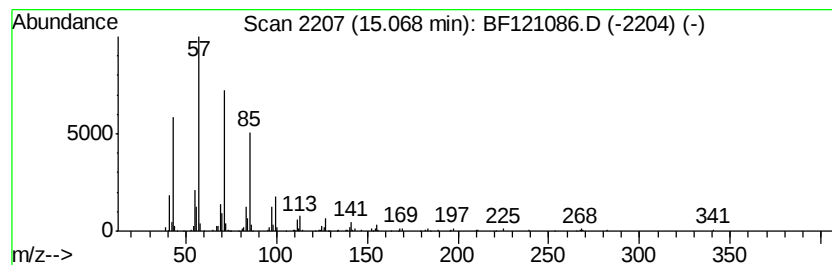
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2-methyloctacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.07	3.48 ng	307606	Perylene-d12		15.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-methyloctacosane	408	C29H60	1000376-72-8	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Eicosane	282	C20H42	000112-95-8	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Nonadecane	268	C19H40	000629-92-5	90



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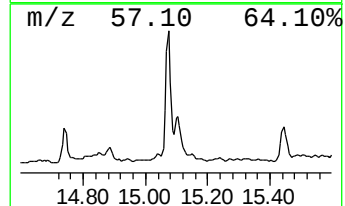
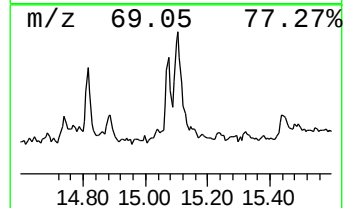
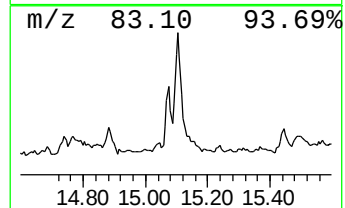
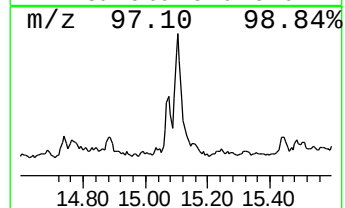
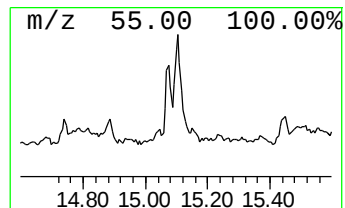
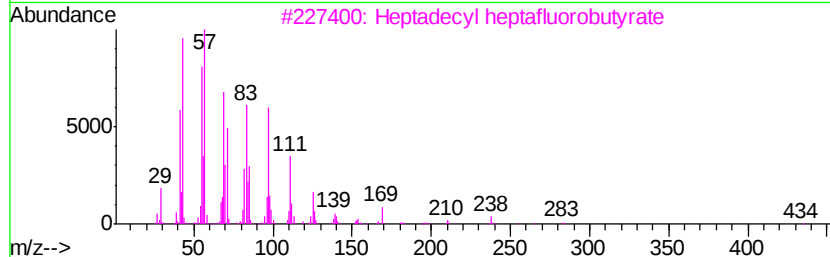
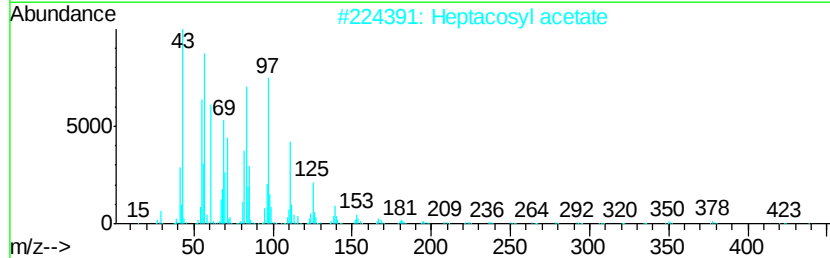
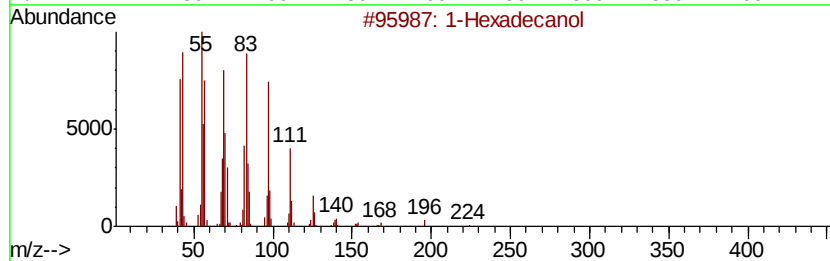
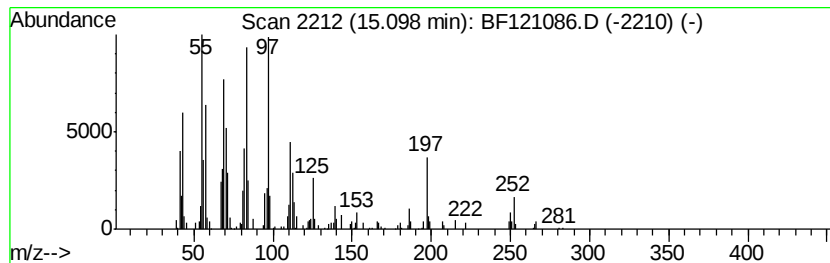
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Hexadecanol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.10	3.05 ng	269833	Perylene-d12	15.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanol	242	C16H34O	036653-82-4	83
2	Heptacosyl acetate	438	C29H58O2	1000351-78-2	83
3	Heptadecyl heptafluorobutrate	452	C21H35F7O2	959085-66-6	83
4	Octacosanol	410	C28H58O	000557-61-9	83
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121086.D
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ClientSampleId :
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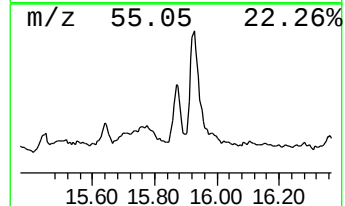
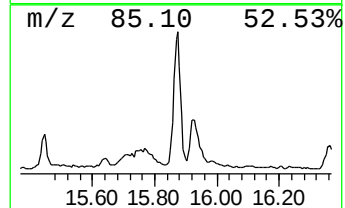
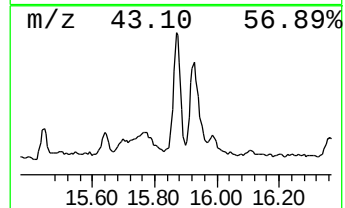
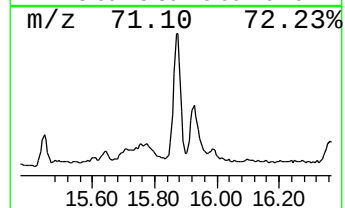
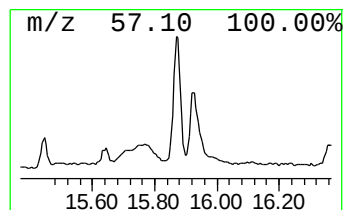
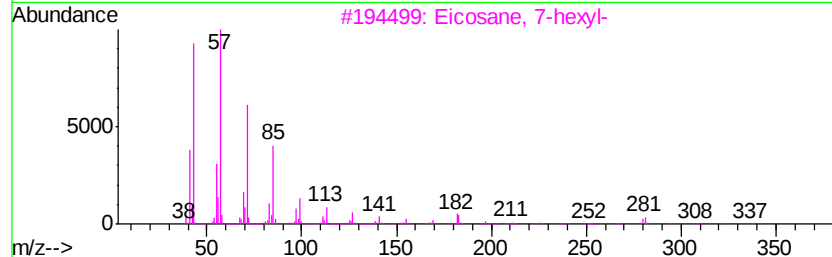
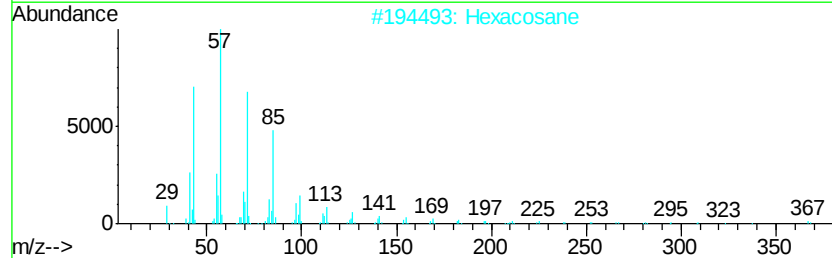
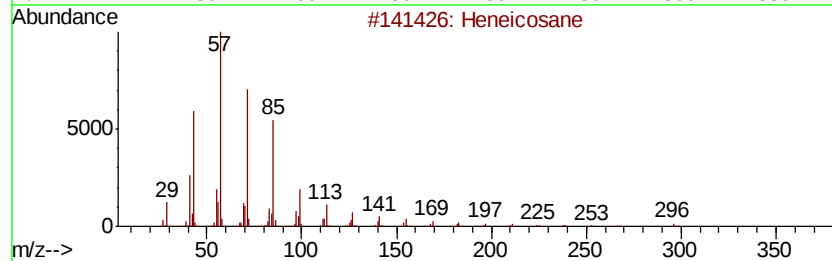
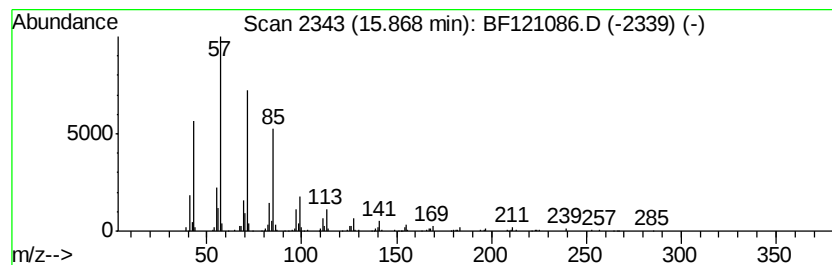
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	5.49 ng	485362	Perylene-d12	15.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Hexacosane		366	C26H54	000630-01-3	91
3	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	90
4	Eicosane		282	C20H42	000112-95-8	90
5	Tetratetracontane		619	C44H90	007098-22-8	87



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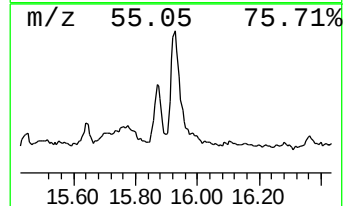
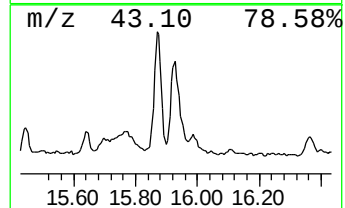
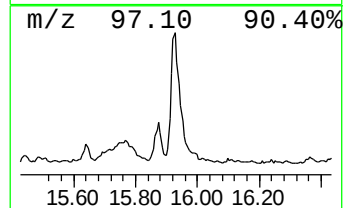
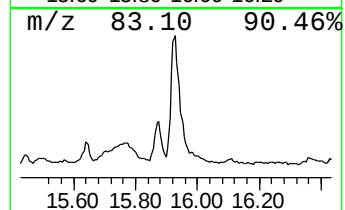
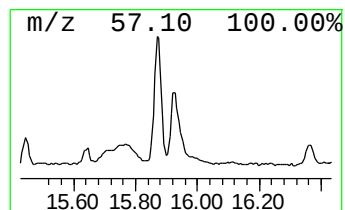
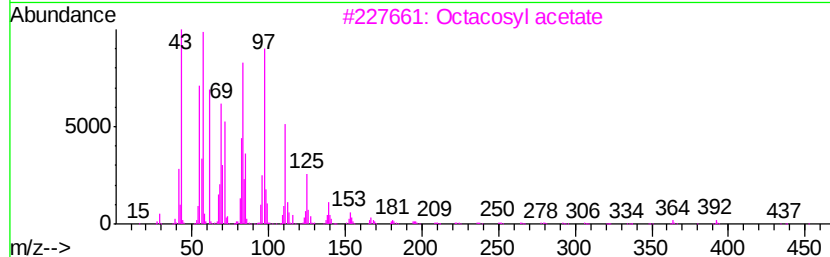
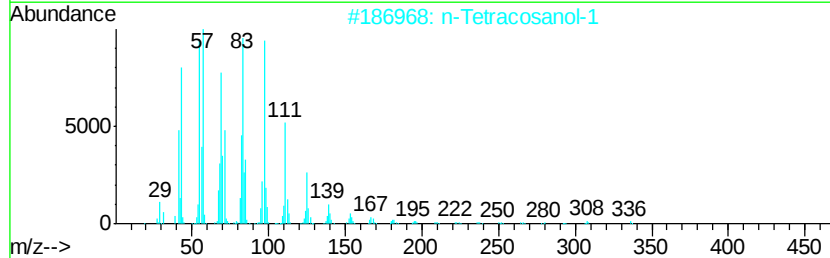
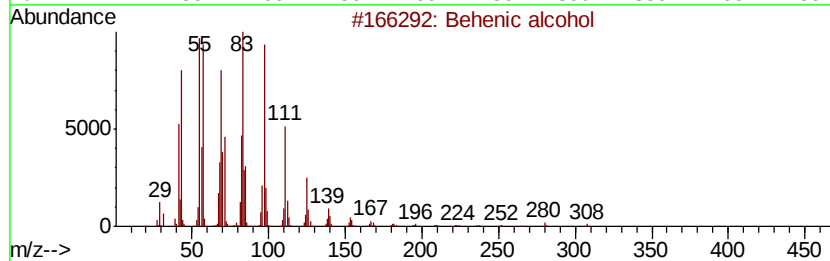
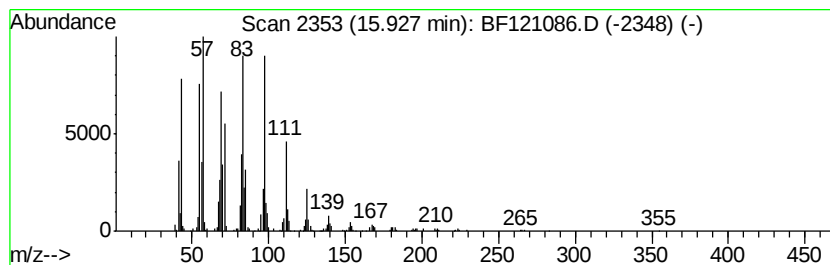
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Behenic alcohol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.93	7.54 ng	666316	Perylene-d12	15.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Behenic alcohol	326	C22H46O	000661-19-8	95
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	95
3	Octacosyl acetate	452	C30H60O2	018206-97-8	95
4	Octacosanol	410	C28H58O	000557-61-9	95
5	1-Heptacosanol	396	C27H56O	002004-39-9	95



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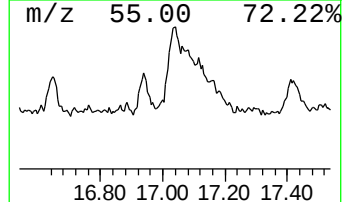
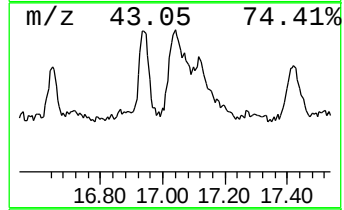
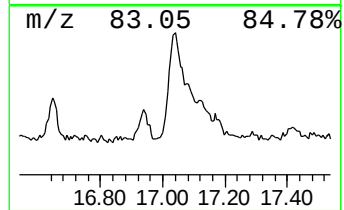
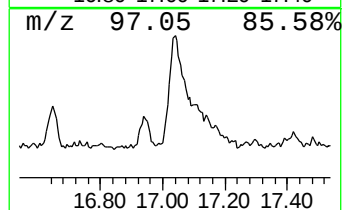
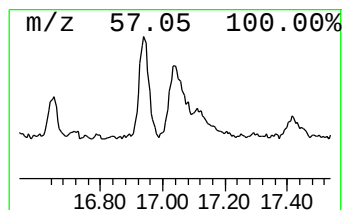
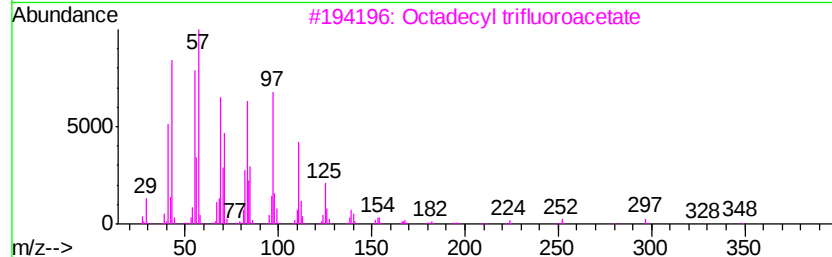
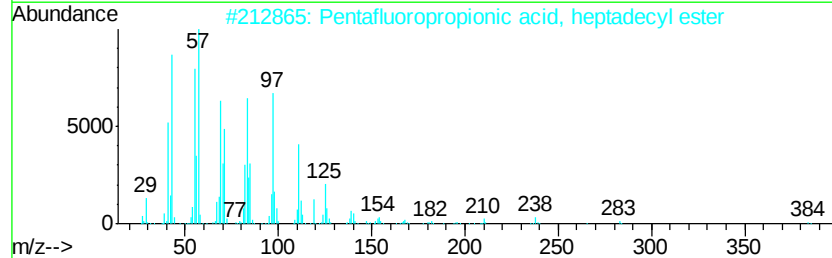
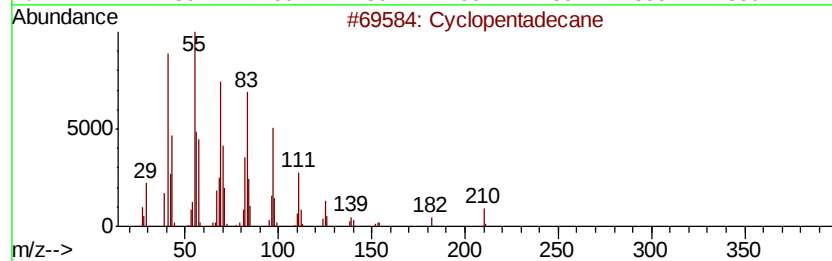
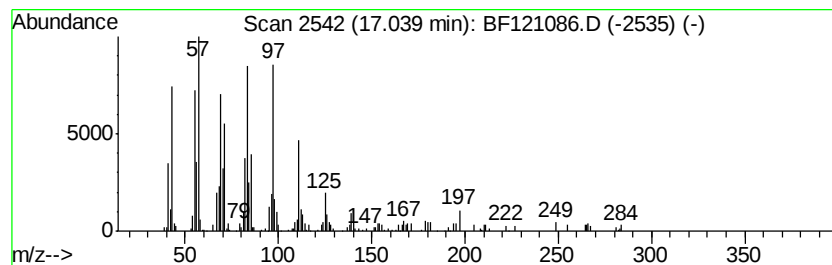
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclopentadecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.04	4.13 ng	364468	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	96
2		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	95
3		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	95
4		Triacetyl acetate	480	C32H64O2	041755-58-2	94
5		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072820\
Data File : BF121086.D
Acq On : 28 Jul 2020 23:58
Operator : JU/CG
Sample : L3435-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

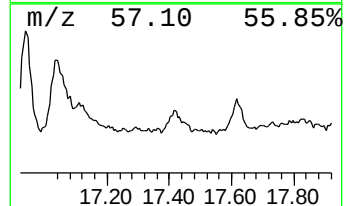
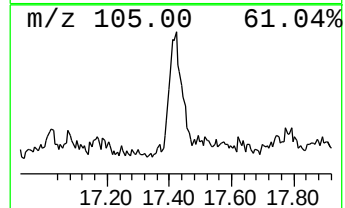
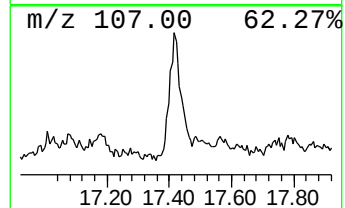
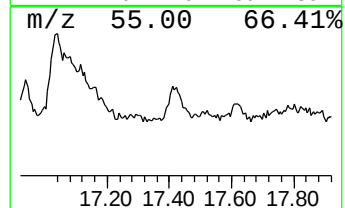
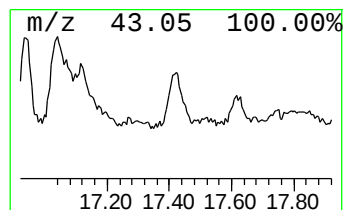
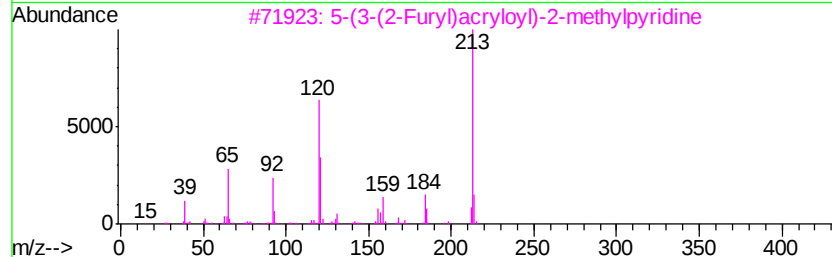
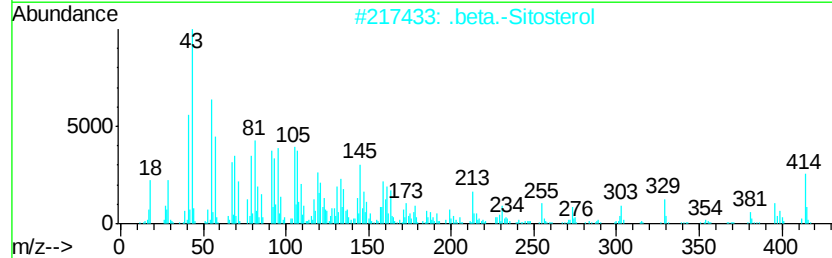
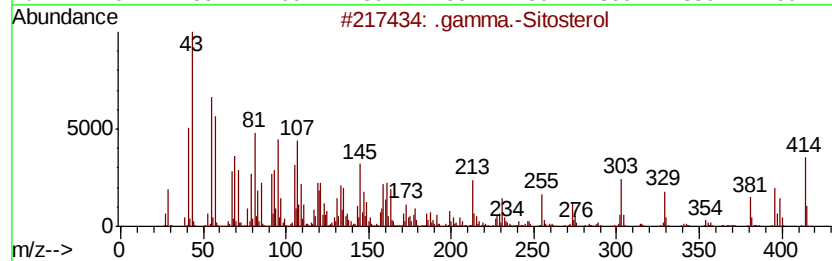
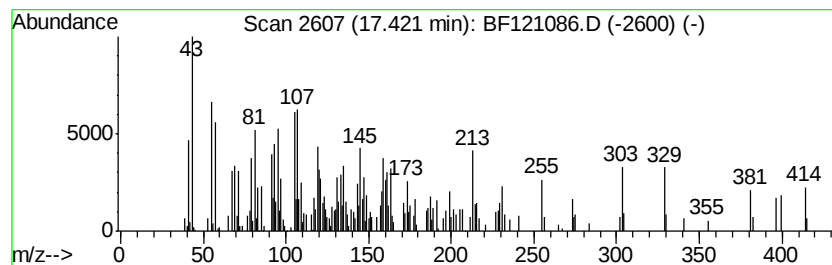
Instrument :
BNA_F
ClientSampleId :
SP-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 .gamma.-Sitosterol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.42	3.79 ng	334432	Perylene-d12		15.42	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		.gamma.-Sitosterol	414	C29H50O	000083-47-6	99
2		.beta.-Sitosterol	414	C29H50O	000083-46-5	96
3		5-(3-(2-Furyl)acryloyl)-2-methylpyridine	213	C13H11NO2	091718-59-1	14
4		Cholest-5-ene, 3-ethoxy-, (3.beta.)-	414	C29H50O	000986-19-6	11
5		N-(3-Chloro-4-fluoro-phenyl)-2-(4-fluorophenyl)ethanamine	329	C14H10Cl2FNOS	1000295-12-7	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072820\
 Data File : BF121086.D
 Acq On : 28 Jul 2020 23:58
 Operator : JU/CG
 Sample : L3435-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF071620.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
Dodecanoic acid	10.06	5.6 ng		367982	3	9.85	1307720	20.0
unknown10.52	10.52	3.4 ng		223999	3	9.85	1307720	20.0
n-Hexadecanoic acid	11.86	4.3 ng		335444	4	11.33	1541360	20.0
2-Phenanthrenol, ...	13.34	2.5 ng		219016	5	13.97	1734960	20.0
5-Eicosene, (E)-	13.82	19.4 ng		1681870	5	13.97	1734960	20.0
unknown14.40	14.40	3.3 ng		288011	5	13.97	1734960	20.0
5-Octadecene, (E)-	14.44	5.6 ng		486086	5	13.97	1734960	20.0
2-methyloctacosane	15.07	3.5 ng		307606	6	15.42	1766710	20.0
1-Hexadecanol	15.10	3.0 ng		269833	6	15.42	1766710	20.0
Heneicosane	15.87	5.5 ng		485362	6	15.42	1766710	20.0
Behenic alcohol	15.93	7.5 ng		666316	6	15.42	1766710	20.0
Cyclopentadecane	17.04	4.1 ng		364468	6	15.42	1766710	20.0
.gamma.-Sitosterol	17.42	3.8 ng		334432	6	15.42	1766710	20.0