

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF081919\
 Data File : BF116412.D
 Acq On : 20 Aug 2019 2:26
 Operator : HP/JU
 Sample : K4223-21
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-11

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-BF073019.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	5.434	566	569	592	rBV	1525703	1876224	66.24%	4.878%
2	6.451	739	742	756	rBV	1815764	2051913	72.44%	5.335%
3	6.581	760	764	773	rVB	2174522	2080723	73.46%	5.410%
4	6.804	799	802	812	rBV	681307	605466	21.37%	1.574%
5	6.957	825	828	846	rVB	1747248	1584089	55.92%	4.119%
6	7.369	895	898	918	rBV	1385604	1335303	47.14%	3.472%
7	8.081	1016	1019	1033	rBV	836695	817048	28.84%	2.124%
8	9.157	1198	1202	1205	rBV	2313411	2150502	75.92%	5.592%
9	9.269	1217	1221	1224	rBV3	42089	40136	1.42%	0.104%
10	9.463	1251	1254	1256	rBV	36456	34245	1.21%	0.089%
11	9.486	1256	1258	1263	rVV	107265	95006	3.35%	0.247%
12	9.551	1263	1269	1272	rVV	444063	378117	13.35%	0.983%
13	9.586	1272	1275	1279	rVB4	55524	52452	1.85%	0.136%
14	9.745	1297	1302	1304	rBV2	34876	35390	1.25%	0.092%
15	9.833	1314	1317	1320	rBV	1177512	988406	34.89%	2.570%
16	9.904	1327	1329	1332	rVB2	40445	33961	1.20%	0.088%
17	9.951	1332	1337	1341	rBV	200714	180831	6.38%	0.470%
18	9.986	1341	1343	1346	rVB	46925	38106	1.35%	0.099%
19	10.063	1350	1356	1358	rBV6	26074	38757	1.37%	0.101%
20	10.128	1363	1367	1369	rVB	49653	53126	1.88%	0.138%
21	10.498	1428	1430	1434	rVB2	65617	57080	2.02%	0.148%
22	10.557	1434	1440	1443	rBV5	51172	77567	2.74%	0.202%
23	10.604	1443	1448	1449	rVV3	73237	102403	3.62%	0.266%
24	10.628	1449	1452	1460	rVV	1676852	1668127	58.89%	4.337%
25	10.686	1460	1462	1466	rVB	41904	41971	1.48%	0.109%
26	10.739	1468	1471	1478	rBV3	58983	66241	2.34%	0.172%
27	10.828	1482	1486	1489	rBV4	21598	31445	1.11%	0.082%
28	10.928	1500	1503	1511	rBV	42054	51731	1.83%	0.135%
29	10.992	1511	1514	1519	rBV4	50379	96116	3.39%	0.250%
30	11.033	1519	1521	1525	rVB	30143	32519	1.15%	0.085%
31	11.322	1564	1570	1573	rVV	1181899	1094764	38.65%	2.846%
32	11.369	1576	1578	1580	rVB	129680	106194	3.75%	0.276%
33	11.451	1590	1592	1596	rVB	90411	84239	2.97%	0.219%
34	11.528	1602	1605	1606	rBV	33945	35159	1.24%	0.091%

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35	11.698	1631	1634	1638	rVB2	45135	56904	2.01%	0.148%
36	11.798	1644	1651	1654	rBV2	114722	187193	6.61%	0.487%
37	11.851	1655	1660	1672	rVV	939068	1124797	39.71%	2.925%
38	12.004	1681	1686	1691	rBV2	213321	253243	8.94%	0.658%
39	12.098	1698	1702	1704	rBV4	45213	52656	1.86%	0.137%
40	12.198	1716	1719	1723	rVB4	46261	48000	1.69%	0.125%
41	12.251	1726	1728	1732	rVV3	55384	60234	2.13%	0.157%
42	12.386	1747	1751	1752	rBV2	48820	54171	1.91%	0.141%
43	12.404	1752	1754	1757	rVB3	86913	77254	2.73%	0.201%
44	12.451	1759	1762	1768	rBV	337902	390216	13.78%	1.015%
45	12.557	1773	1780	1789	rBV5	750726	1763423	62.25%	4.585%
46	12.627	1789	1792	1797	rBV	214975	228246	8.06%	0.593%
47	12.722	1805	1808	1811	rBV2	104568	100460	3.55%	0.261%
48	12.769	1813	1816	1821	rVB	184506	186759	6.59%	0.486%
49	12.898	1834	1838	1842	rBV	2231383	1909992	67.43%	4.966%
50	13.098	1870	1872	1874	rBV2	42743	36215	1.28%	0.094%
51	13.122	1874	1876	1881	rVB3	66837	62496	2.21%	0.162%
52	13.292	1901	1905	1907	rBV3	106983	130013	4.59%	0.338%
53	13.327	1908	1911	1912	rBV	215942	199242	7.03%	0.518%
54	13.392	1920	1922	1924	rBV	51348	46374	1.64%	0.121%
55	13.416	1924	1926	1930	rVB3	94685	85525	3.02%	0.222%
56	13.580	1952	1954	1958	rBV3	40336	39124	1.38%	0.102%
57	13.786	1985	1989	1995	rBV	430907	482417	17.03%	1.254%
58	13.874	2001	2004	2008	rBV	96082	103243	3.64%	0.268%
59	13.963	2016	2019	2023	rBV	768767	768594	27.13%	1.998%
60	14.104	2041	2043	2045	rBV	64509	48430	1.71%	0.126%
61	14.410	2091	2095	2102	rBV2	550050	760273	26.84%	1.977%
62	14.704	2143	2145	2148	rVB	110977	98484	3.48%	0.256%
63	14.780	2155	2158	2161	rVB	244587	237142	8.37%	0.617%
64	14.927	2180	2183	2187	rBV3	76714	97960	3.46%	0.255%
65	15.033	2197	2201	2204	rBV	1491959	1574783	55.59%	4.095%
66	15.063	2204	2206	2212	rVB	179251	198349	7.00%	0.516%
67	15.286	2242	2244	2251	rVB2	87974	117607	4.15%	0.306%
68	15.392	2260	2262	2264	rBV	77937	89572	3.16%	0.233%
69	15.421	2264	2267	2276	rVB	608585	782568	27.63%	2.035%
70	15.815	2329	2334	2339	rVB	1108443	1553263	54.83%	4.039%
71	15.868	2339	2343	2351	rVB2	165739	267883	9.46%	0.697%

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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

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72	16.086	2376	2380	2391	rVB	297934	572182	20.20%	1.488%
73	17.304	2583	2587	2592	rBV3	96504	193133	6.82%	0.502%
74	17.380	2593	2600	2613	rVV6	798674	2832618	100.00%	7.365%
75	17.821	2669	2675	2685	rBV	337384	918353	32.42%	2.388%
76	18.151	2724	2731	2736	rBV2	436469	1188618	41.96%	3.091%
77	18.374	2765	2769	2780	rVB5	178291	466678	16.48%	1.213%

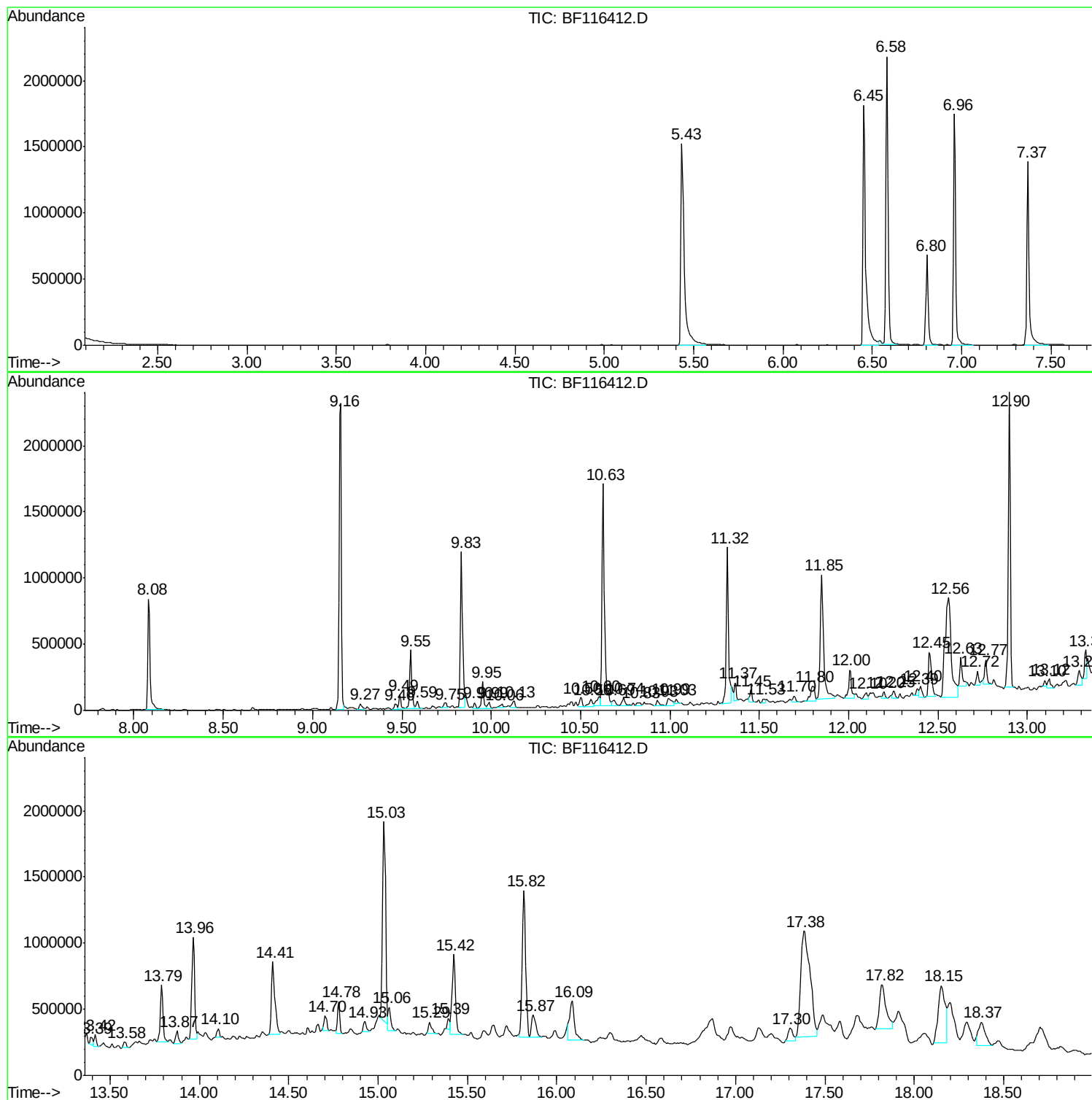
Sum of corrected areas: 38460044

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Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
SP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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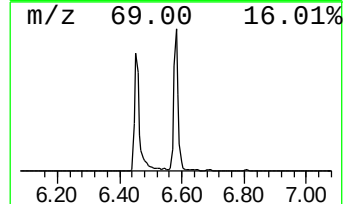
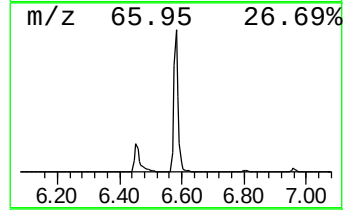
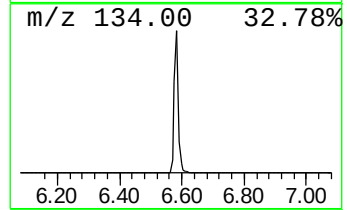
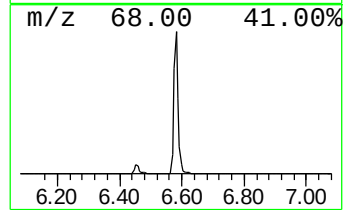
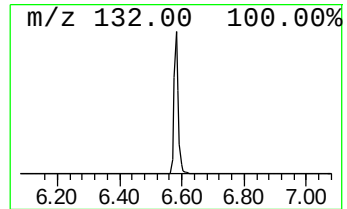
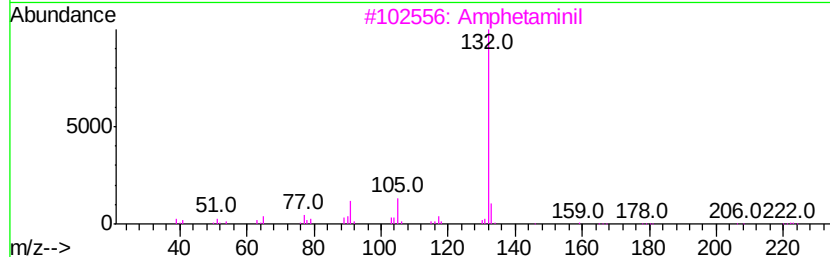
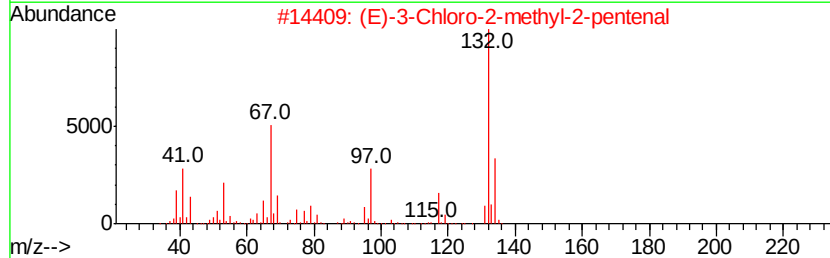
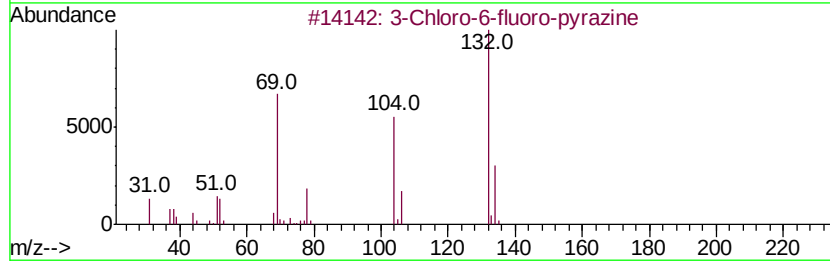
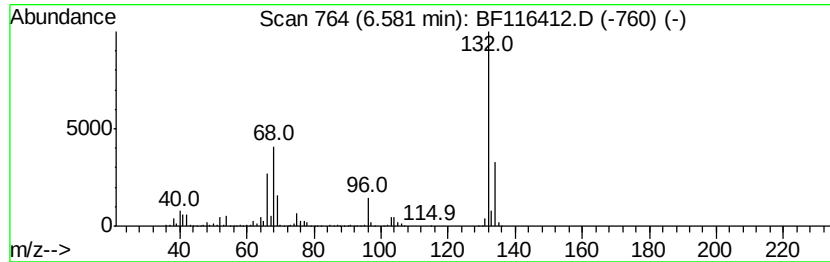
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	68.73 ng	2080720	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Xenon	132	Xe	007440-63-3	9



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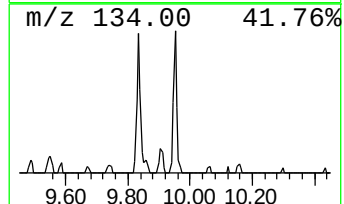
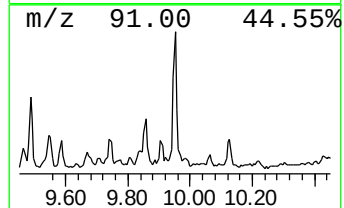
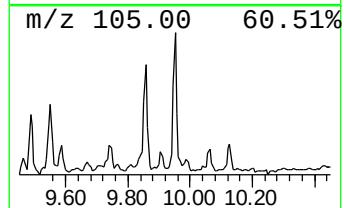
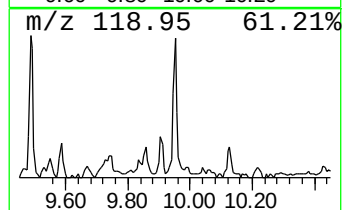
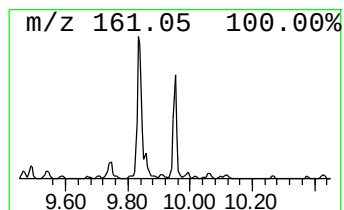
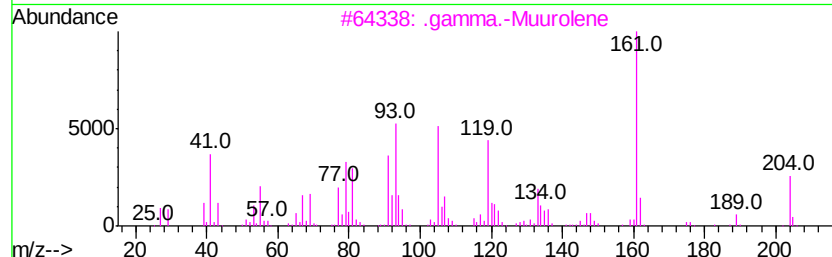
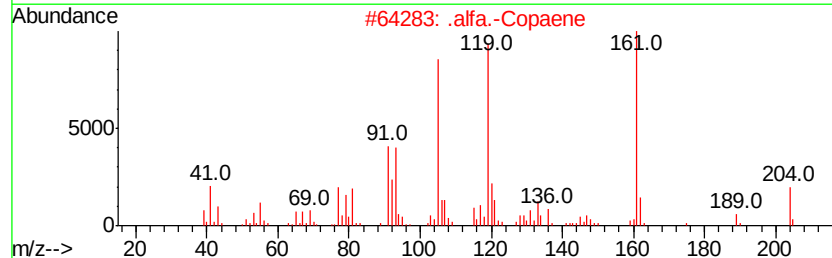
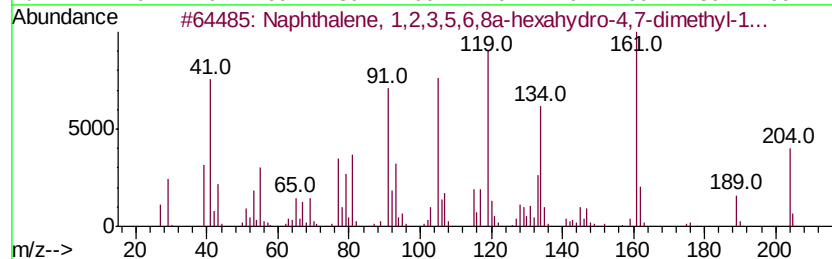
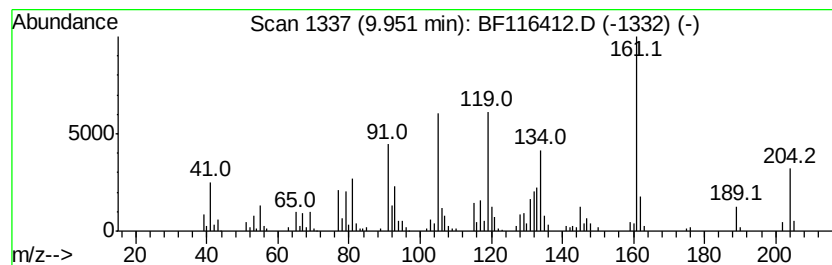
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, 1,2,3,5,6,8a-h... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.95	3.66 ng	180831	Acenaphthene-d10	9.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,5,6,8a-hexahy...	204	C15H24	000483-76-1	95
2		.alfa.-Copaene	204	C15H24	1000360-33-0	91
3		.gamma.-Muurolene	204	C15H24	030021-74-0	86
4		Copaene	204	C15H24	003856-25-5	83
5		Epizonarene	204	C15H24	041702-63-0	76



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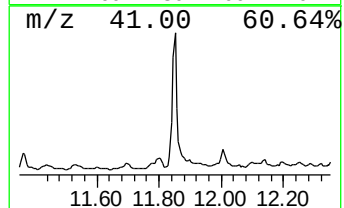
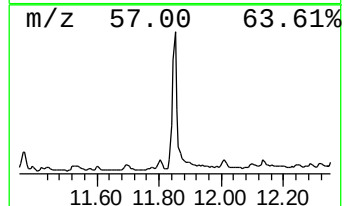
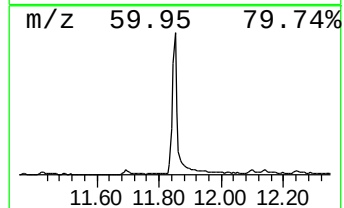
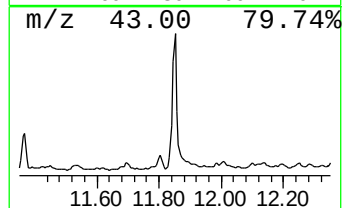
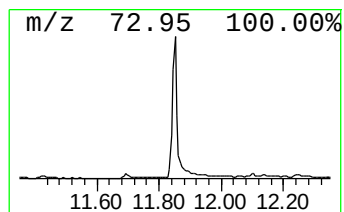
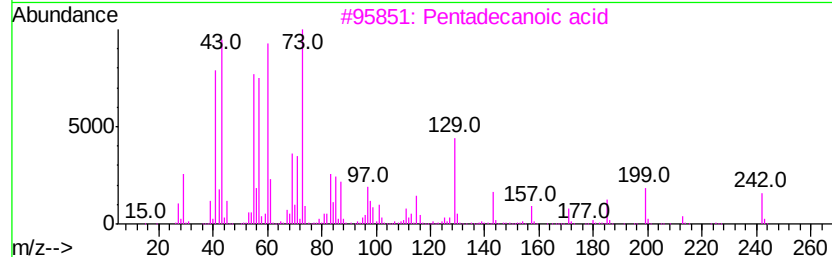
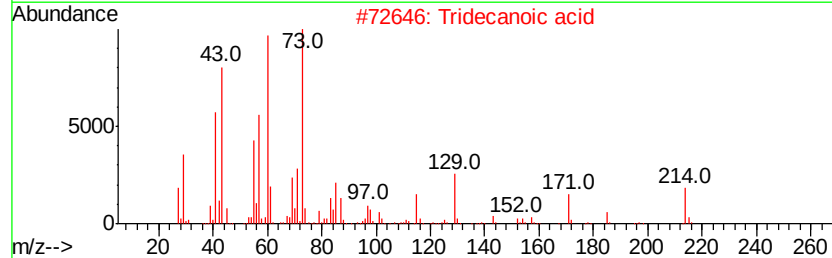
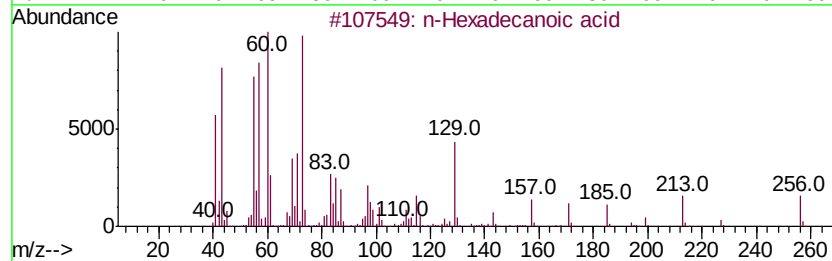
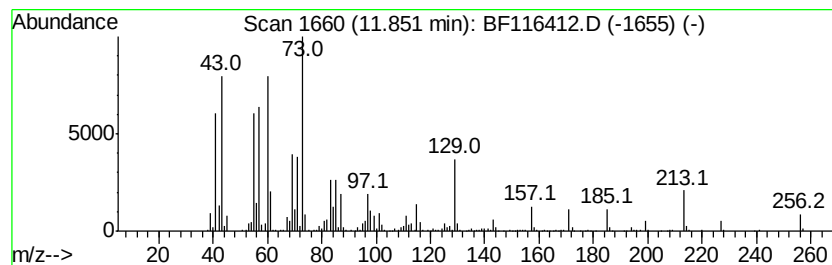
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.85	20.55 ng	1124800	Phenanthrene-d10	11.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tridecanoic acid	214	C13H26O2	000638-53-9	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
5		Oxalic acid, allyl octadecyl ester	382	C23H42O4	1000309-24-5	18



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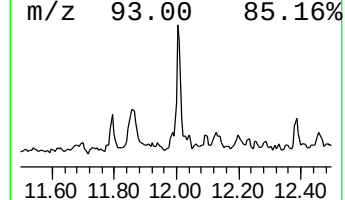
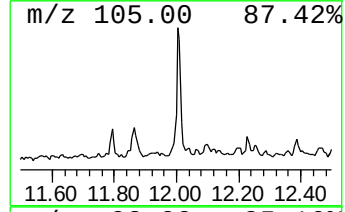
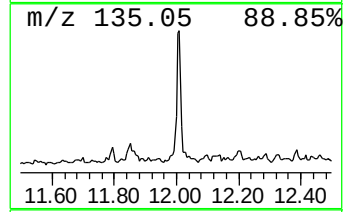
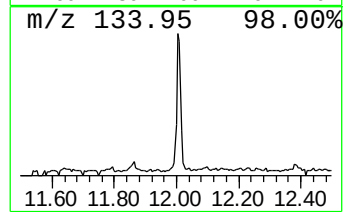
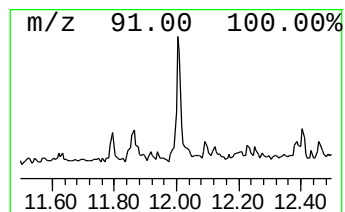
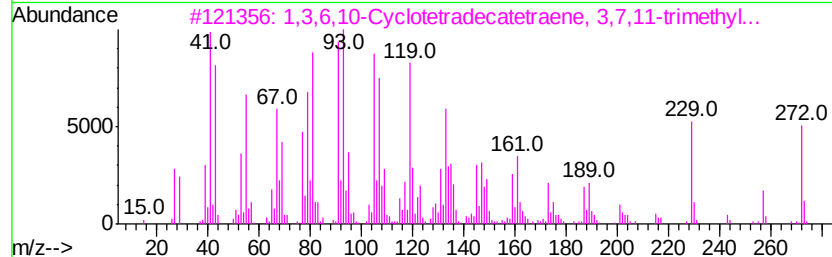
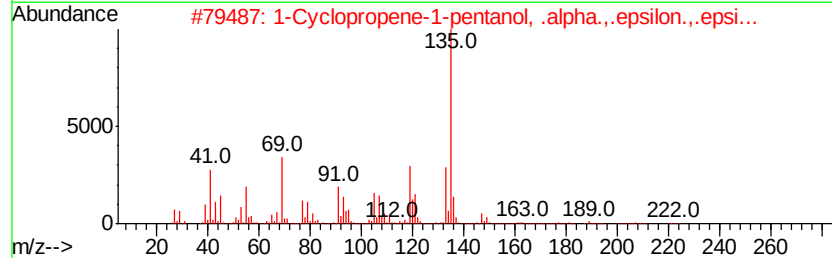
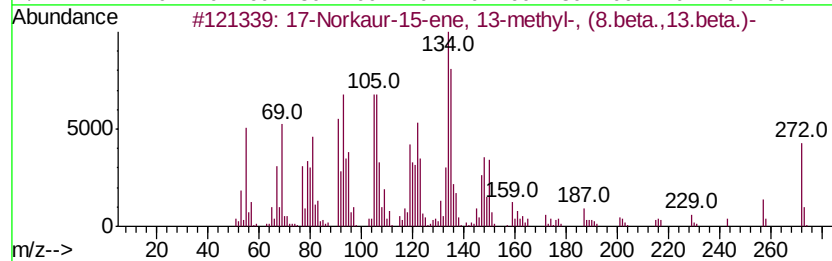
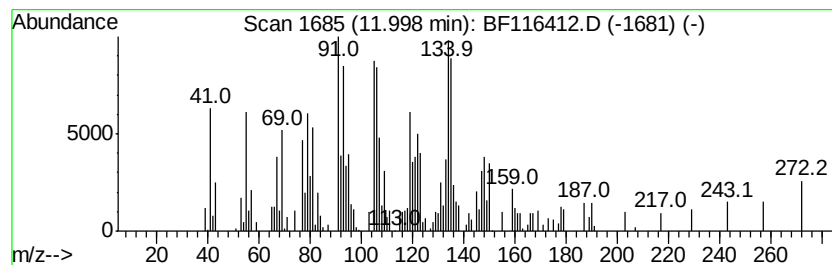
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 17-Norkaur-15-ene, 13-methy... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.00	4.63 ng	253243	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-Norkaur-15-ene, 13-methyl-, (...	272	C20H32	003564-54-3	99
2	1-Cyclopropene-1-pentanol, .alph...	222	C15H26O	090165-06-3	30
3	1,3,6,10-Cyclotetradecatetraene,...	272	C20H32	001898-13-1	25
4	4-Methylformanilide	135	C8H9NO	003085-54-9	25
5	Benzenamine, N,N,3-trimethyl-	135	C9H13N	000121-72-2	22



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
Data File : BF116412.D
Acq On : 20 Aug 2019 2:26
Operator : HP/JU
Sample : K4223-21
Misc :
ALS Vial : 21 Sample Multiplier: 1

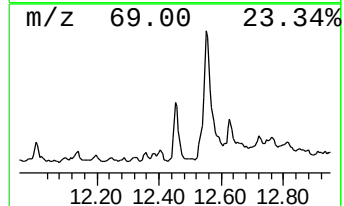
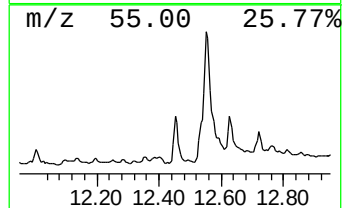
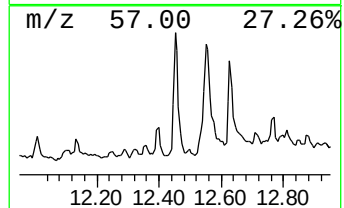
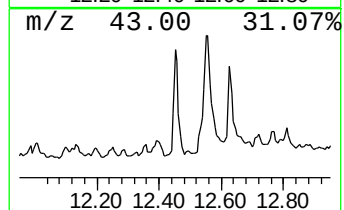
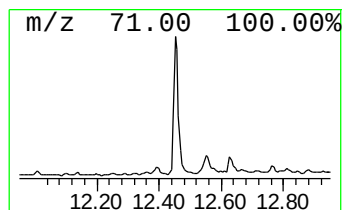
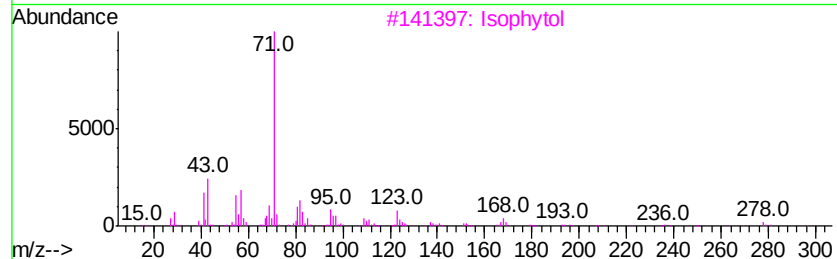
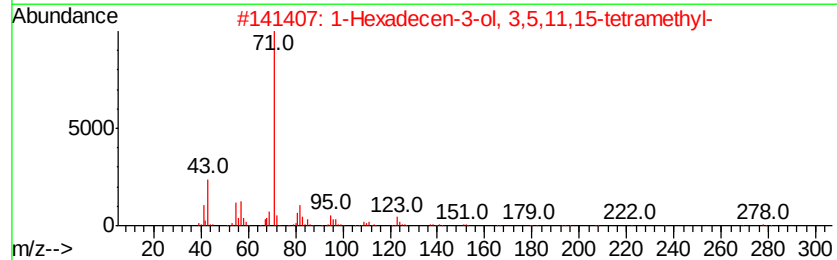
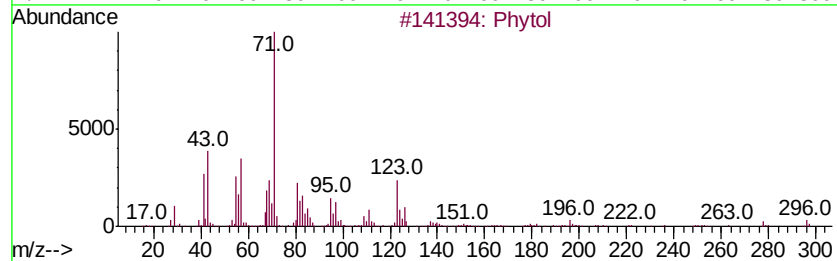
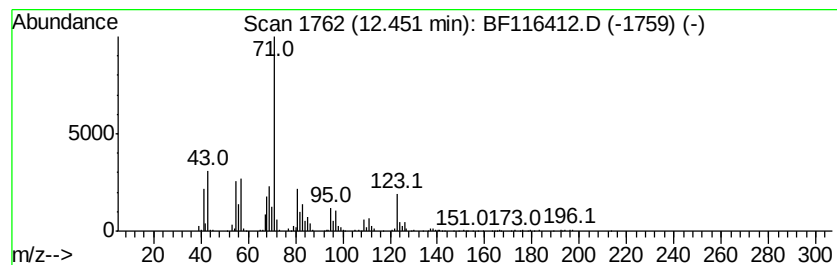
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Phytol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.45	7.13 ng	390216	Phenanthrene-d10		11.32	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phytol		296	C20H40O	000150-86-7	86
2	1-Hexadecen-3-ol, 3,5,11,15-tetr...		296	C20H40O	1000262-55-5	59
3	Isophytol		296	C20H40O	000505-32-8	59
4	Butyric acid, 3-tetradecyl ester		284	C18H36O2	1000280-57-1	52
5	Malonic acid, neopentyl tridecyl...		356	C21H40O4	1000363-93-6	50



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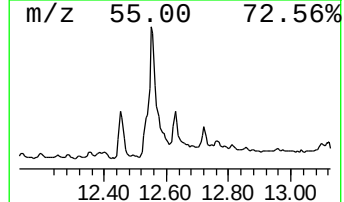
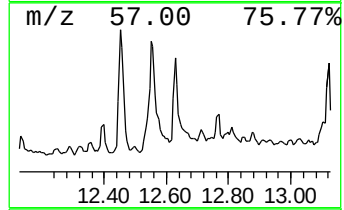
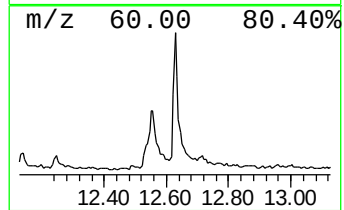
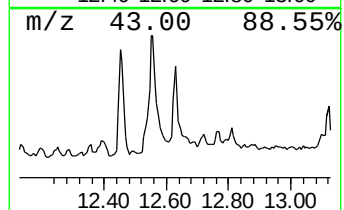
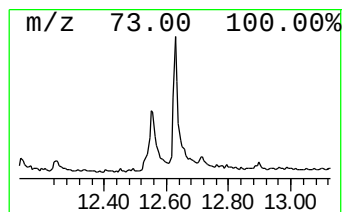
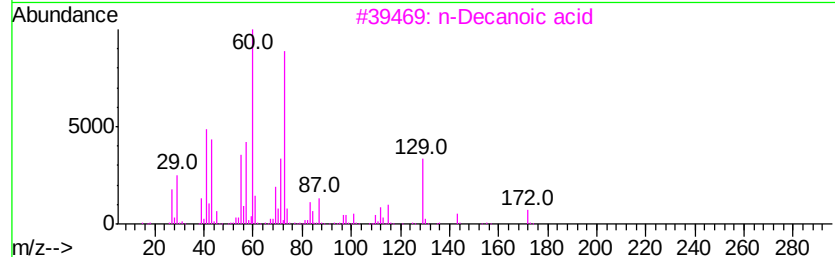
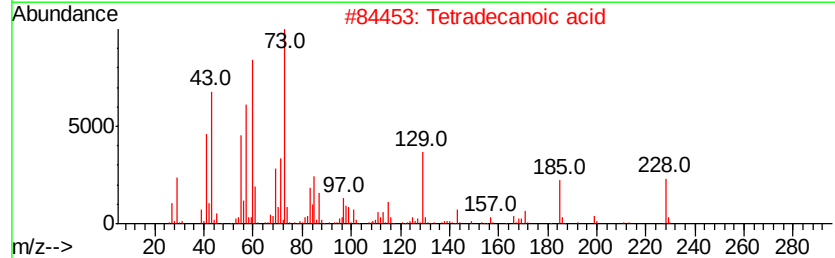
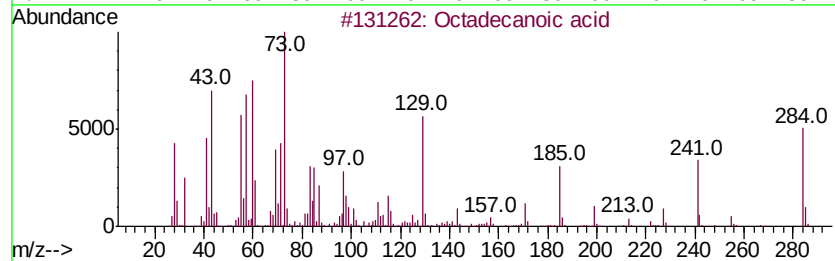
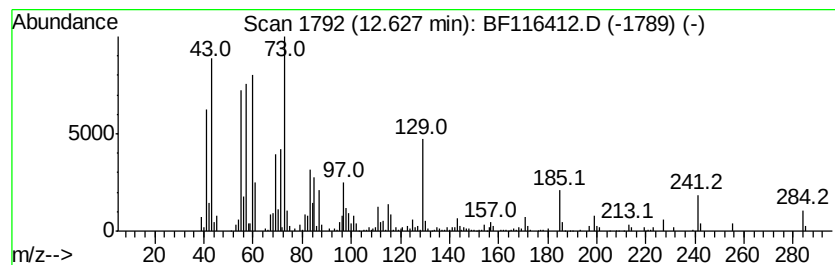
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.63	4.17 ng	228246	Phenanthrene-d10	11.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
3	n-Decanoic acid	172	C10H20O2	000334-48-5	76
4	Tridecanoic acid	214	C13H26O2	000638-53-9	76
5	Octadecanoic acid, 2-(2-hydroxy...	372	C22H44O4	000106-11-6	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
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Sample : K4223-21
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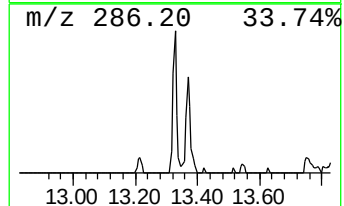
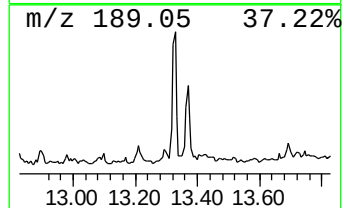
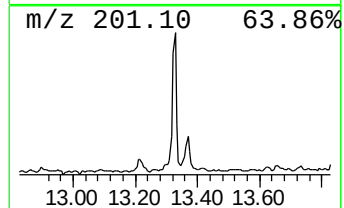
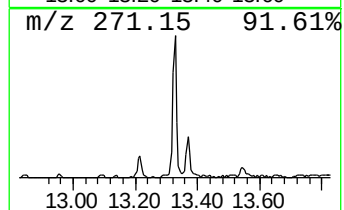
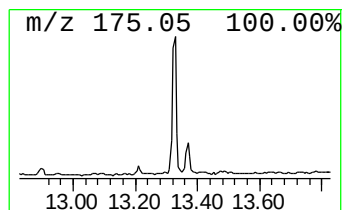
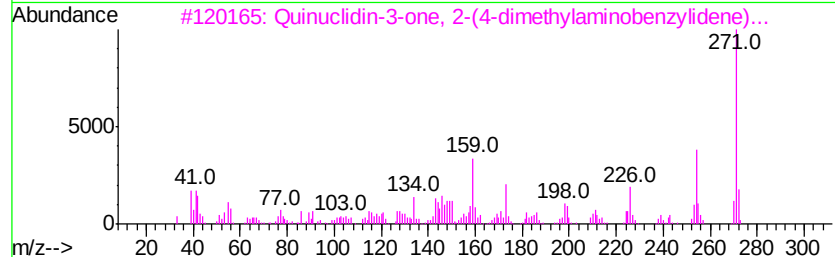
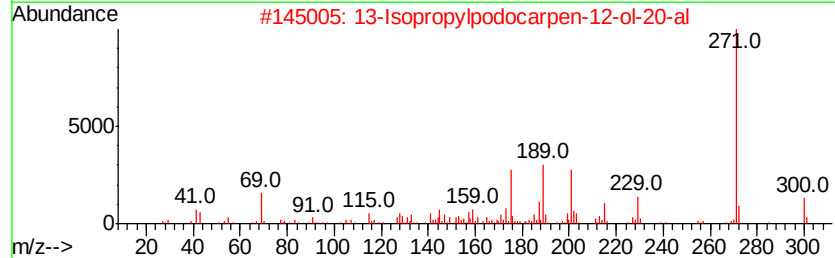
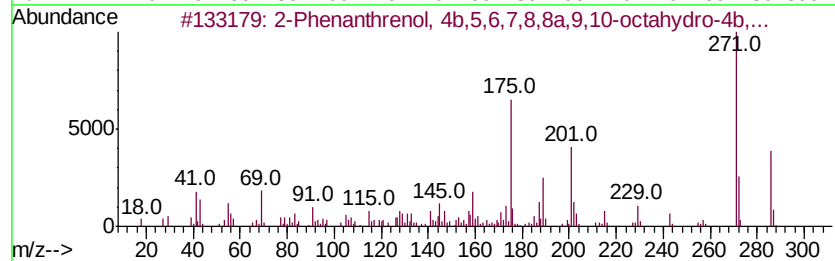
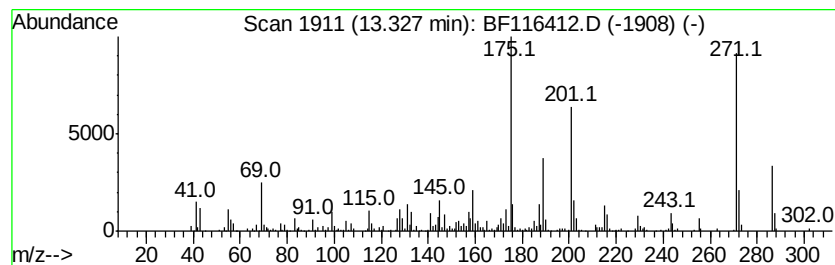
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.33	5.18 ng	199242	Chrysene-d12	13.96		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	93	
2	13-Isopropylpodocarpin-12-ol-20-al	300	C20H28O2	024035-37-8	35	
3	Quinuclidin-3-one, 2-(4-dimethyl...	271	C16H21N3O	1000266-14-6	27	
4	1-Methyl-3-phenyl-4-azafluorenone	271	C19H13NO	069751-55-9	22	
5	Methanone, (2-chloro-4,5-difluor...	287	C14H16ClF2NO	1000277-48-2	22	



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
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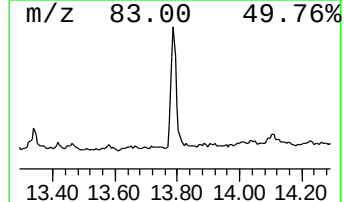
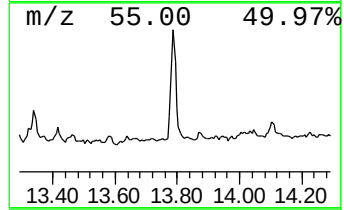
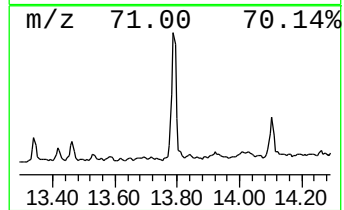
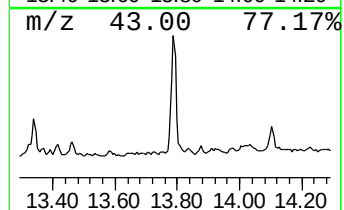
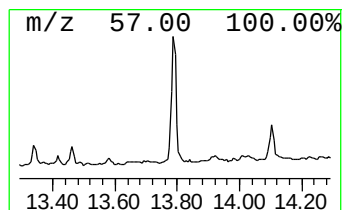
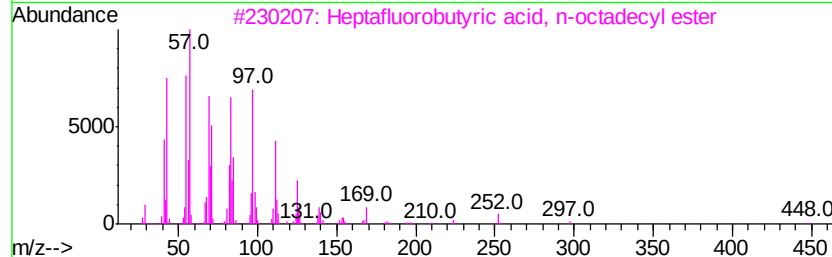
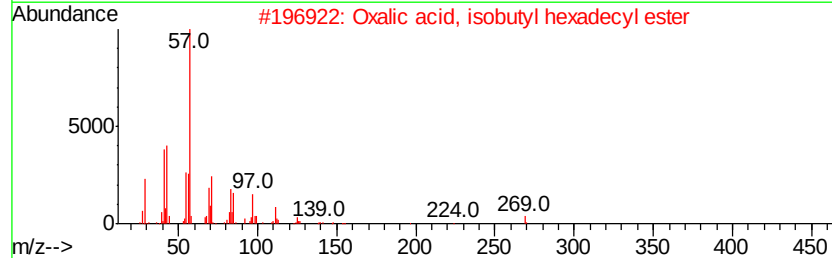
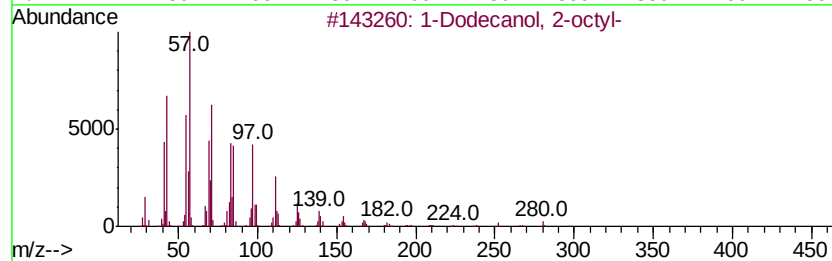
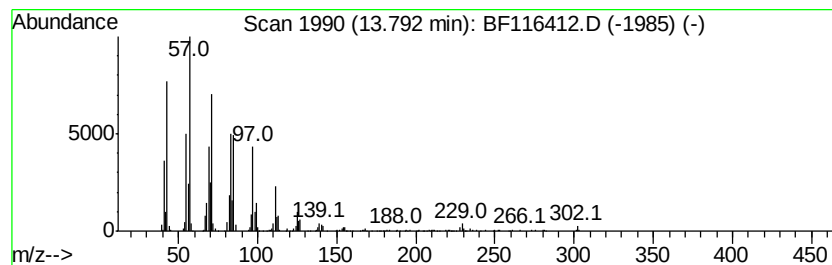
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Dodecanol, 2-octyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	12.55 ng	482417	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	91
2		Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
3		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	76
4		Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	76
5		1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	74



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
Data File : BF116412.D
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Misc :
ALS Vial : 21 Sample Multiplier: 1

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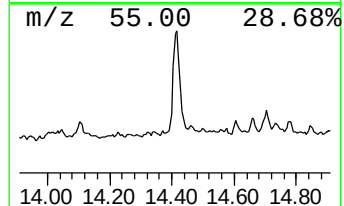
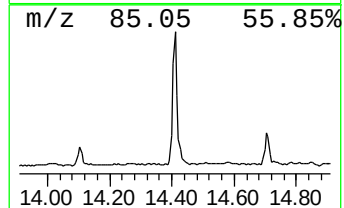
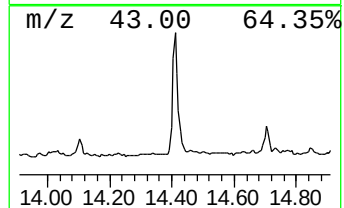
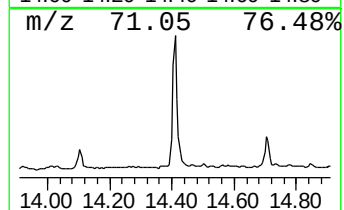
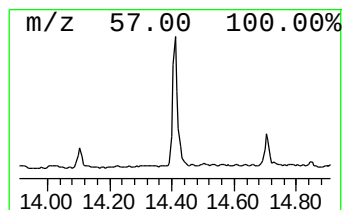
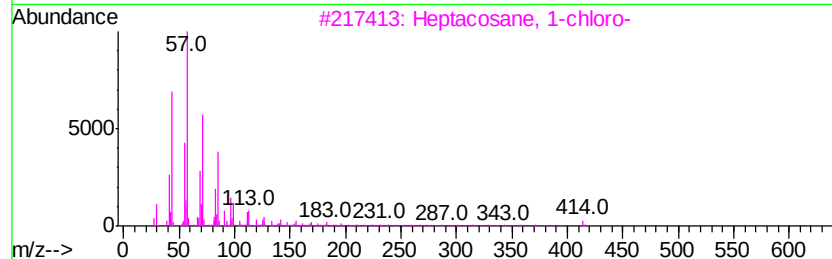
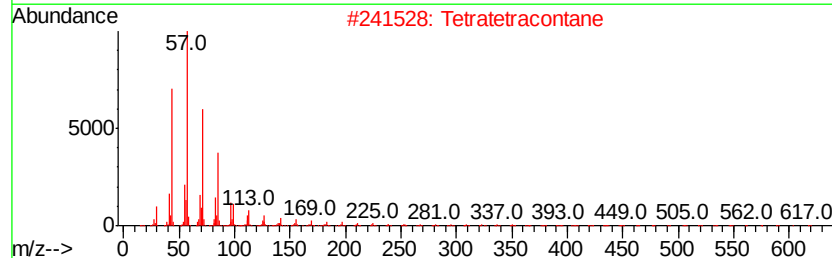
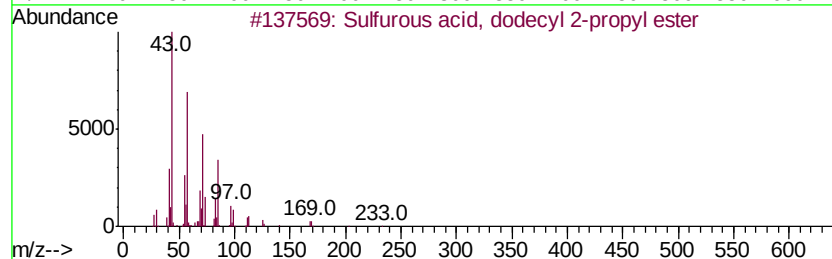
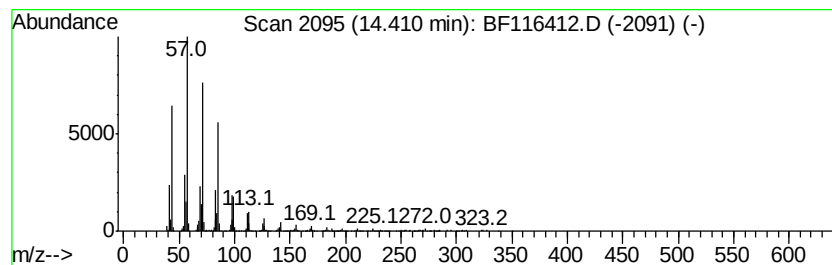
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Sulfurous acid, dodecyl 2-p... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.41	19.78 ng	760273	Chrysene-d12	13.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	91
2		Tetratetracontane	619	C44H90	007098-22-8	91
3		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	91
4		Hexadecane	226	C16H34	000544-76-3	90
5		Hexacosane	366	C26H54	000630-01-3	87



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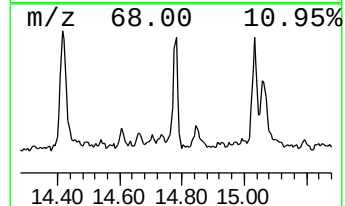
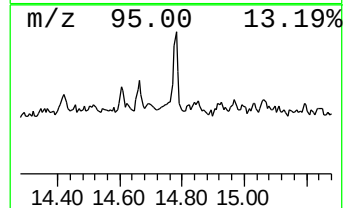
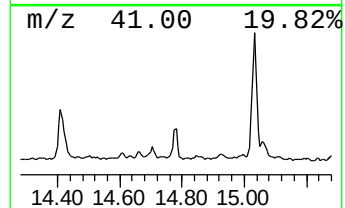
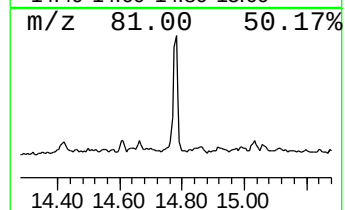
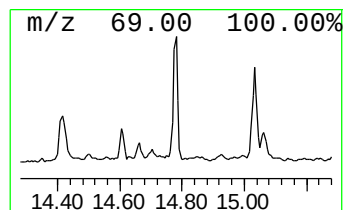
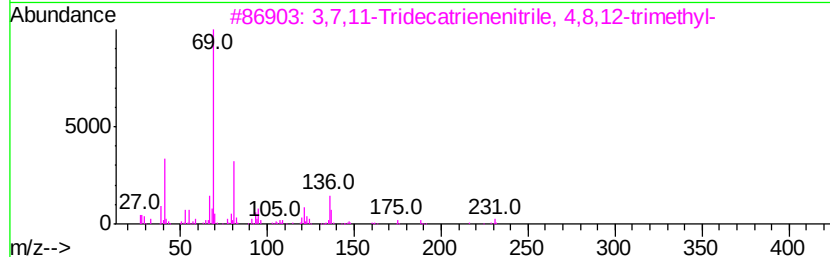
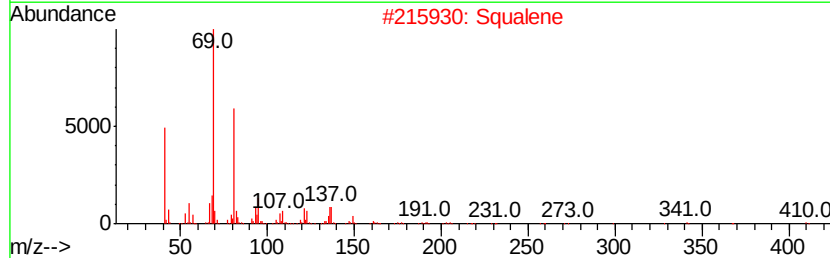
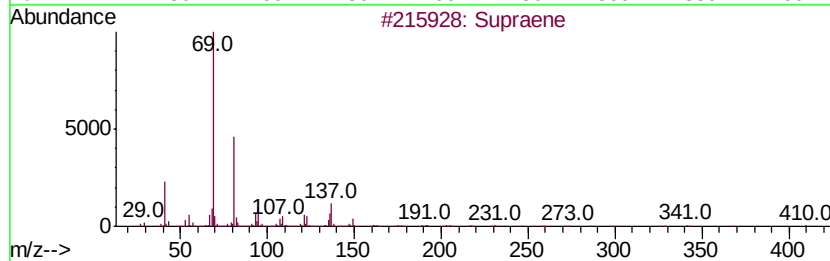
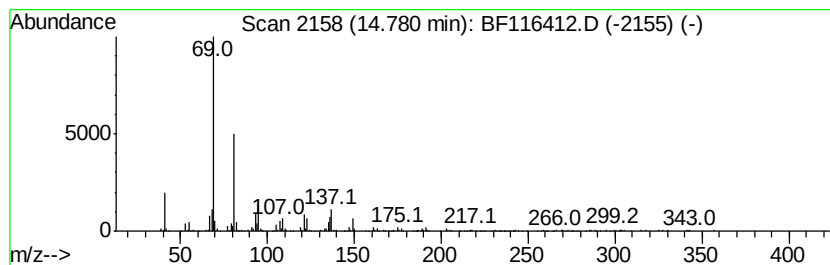
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Squalene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.78	6.06 ng	237142	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Supraene	410	C30H50	007683-64-9	91
2		Squalene	410	C30H50	000111-02-4	91
3		3,7,11-Tridecatrienitrile, 4,8...	231	C16H25N	006006-01-5	72
4		2,6-Octadiene, 4-methyl-	124	C9H16	074498-94-5	50
5		Butanoic acid, 4-cyano-2-ethyl-2...	290	C15H18N2O4	203808-68-8	49



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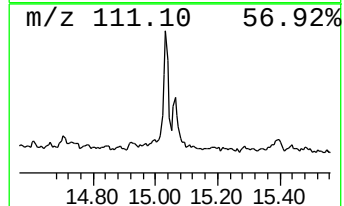
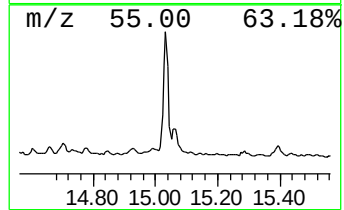
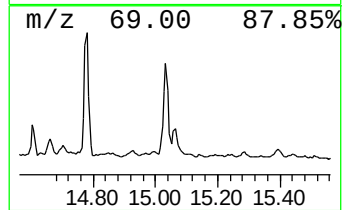
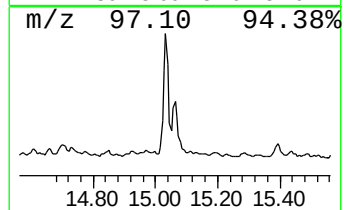
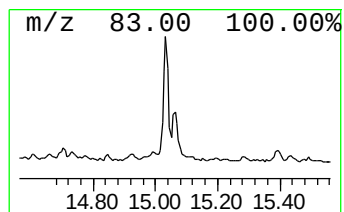
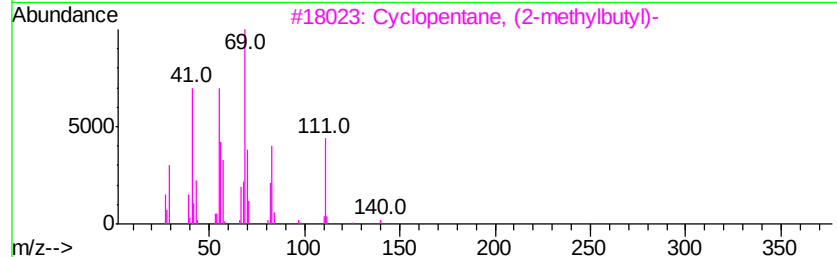
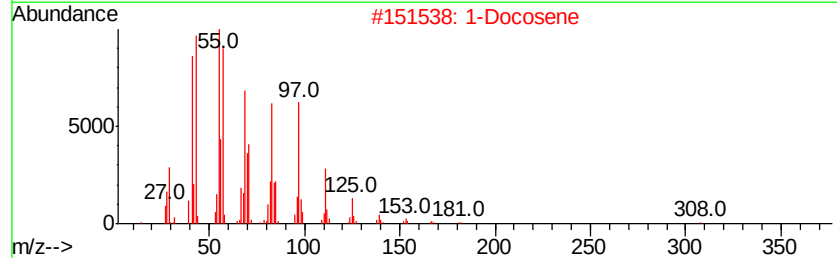
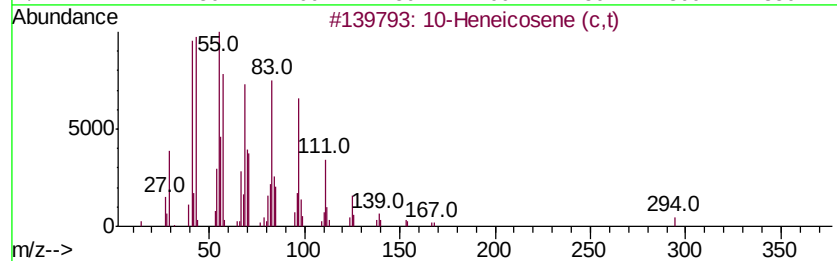
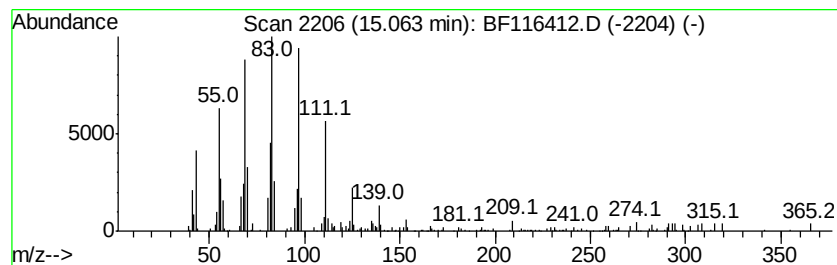
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 10-Heneicosene (c,t) Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	5.07 ng	198349	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10-Heneicosene (c,t)	294	C21H42	095008-11-0	70
2		1-Docosene	308	C22H44	001599-67-3	49
3		Cyclopentane, (2-methylbutyl)-	140	C10H20	053366-38-4	42
4		Cyclooctane, cyclohexyl-	194	C14H26	092369-78-3	41
5		Cyclohexane, (1,2-dimethylbutyl)-	168	C12H24	061142-37-8	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
Data File : BF116412.D
Acq On : 20 Aug 2019 2:26
Operator : HP/JU
Sample : K4223-21
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-11

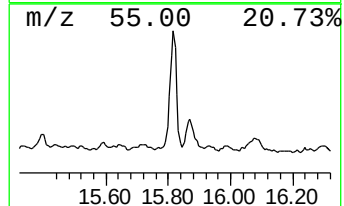
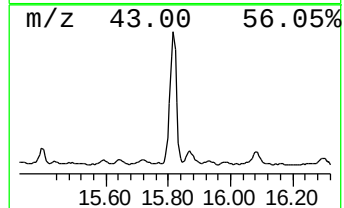
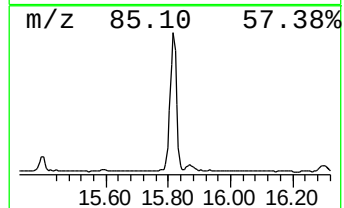
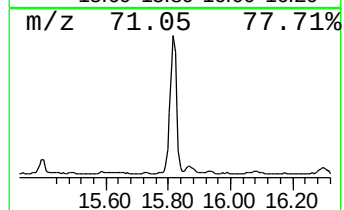
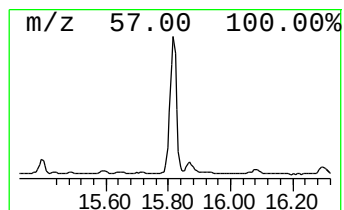
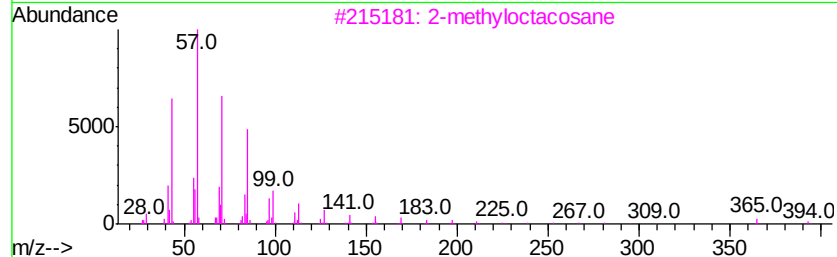
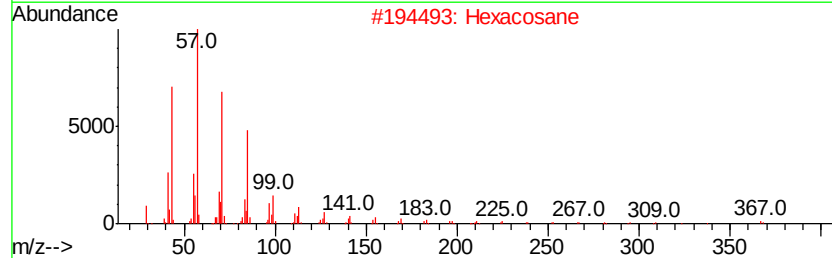
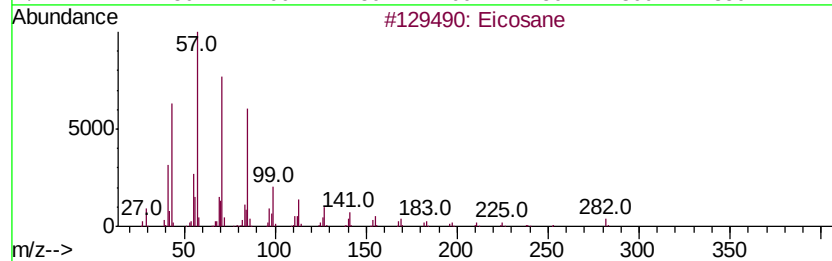
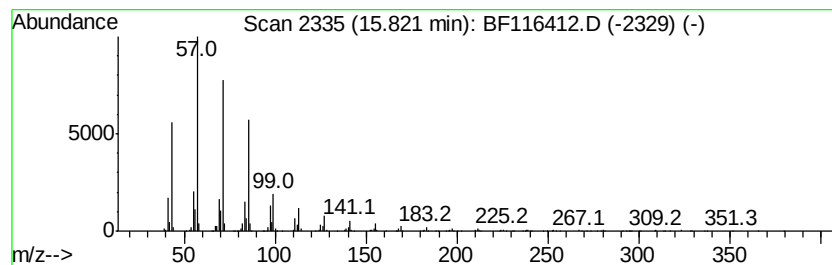
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Eicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	39.70 ng	1553260	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	95
2		Hexacosane	366	C26H54	000630-01-3	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Heneicosane	296	C21H44	000629-94-7	91
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87



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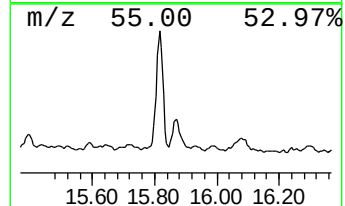
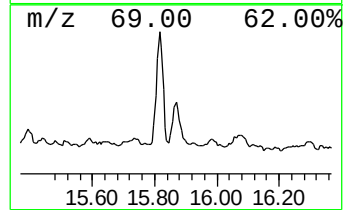
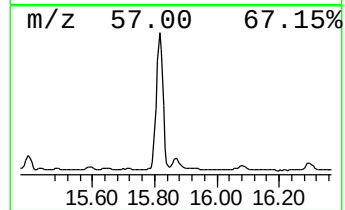
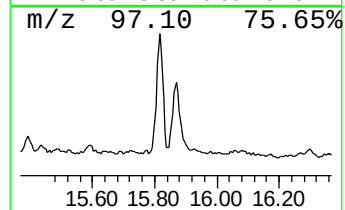
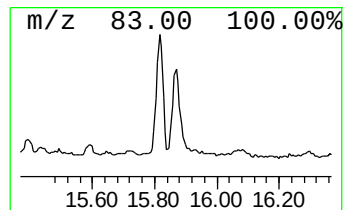
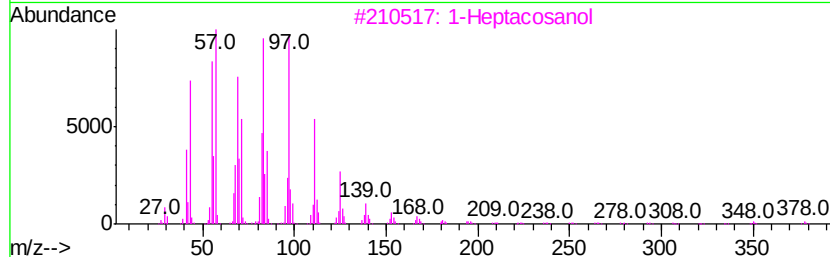
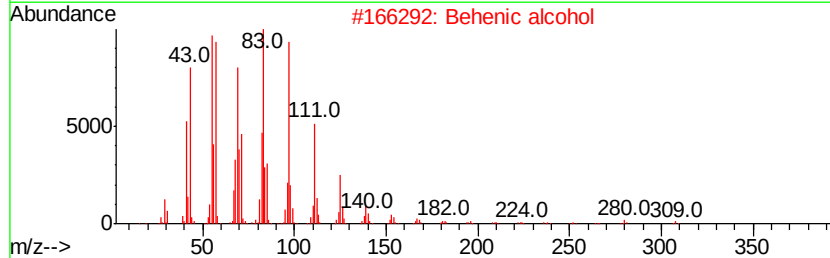
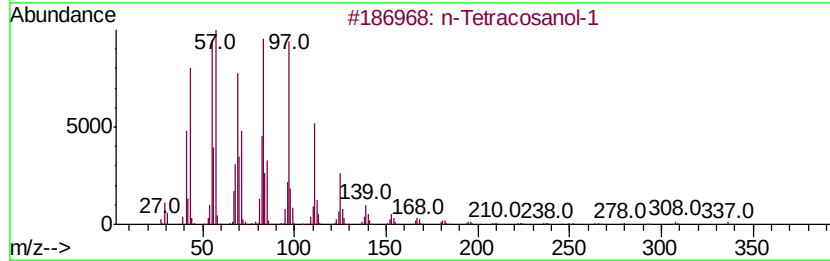
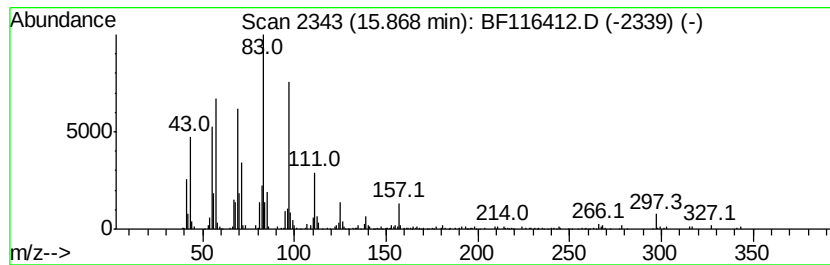
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 n-Tetracosanol-1 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.87	6.85 ng	267883	Perylene-d12	15.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Tetracosanol-1	354	C24H50O	000506-51-4	60
2	Behenic alcohol	326	C22H46O	000661-19-8	53
3	1-Heptacosanol	396	C27H56O	002004-39-9	53
4	Cycloheptane, methyl-	112	C8H16	004126-78-7	50
5	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	41



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
Data File : BF116412.D
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Sample : K4223-21
Misc :
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Instrument :
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ClientSampleId :
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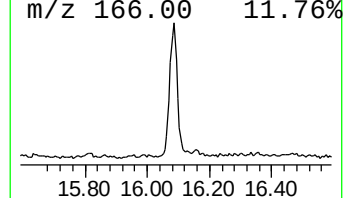
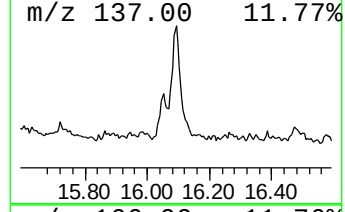
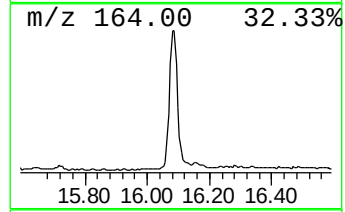
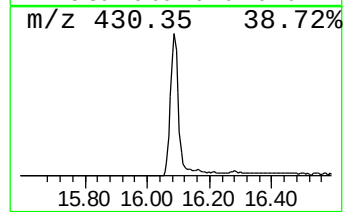
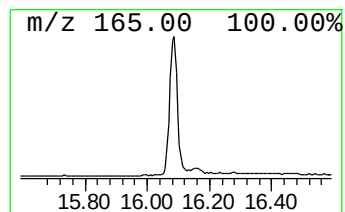
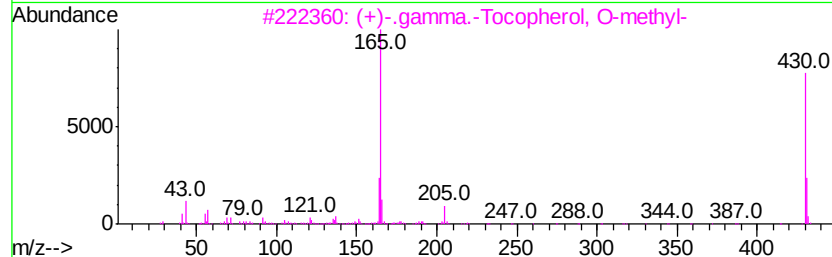
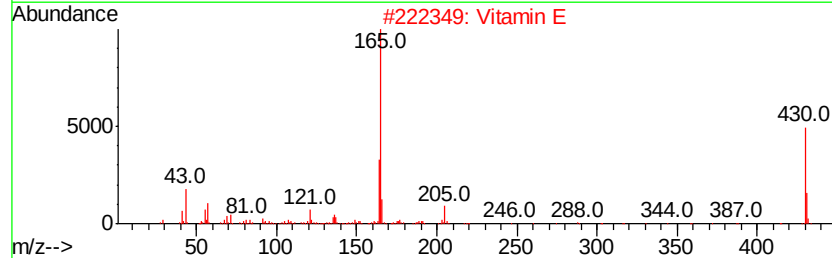
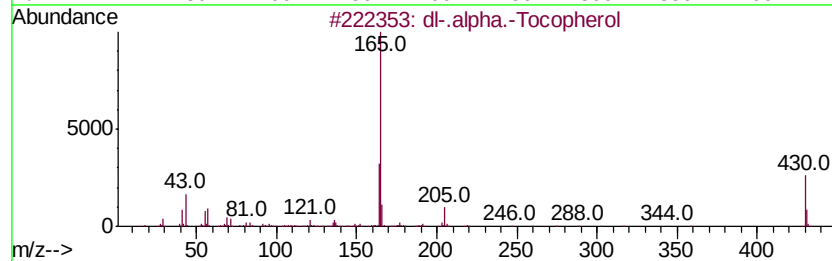
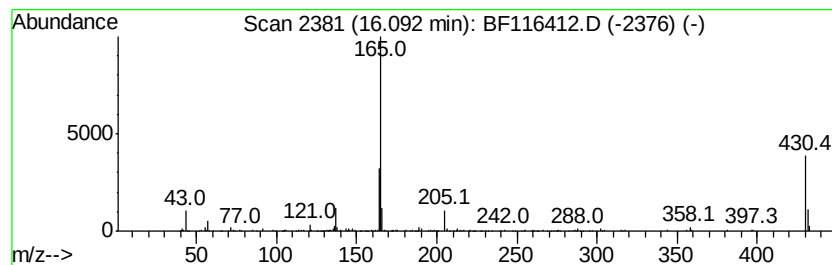
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 dl-.alpha.-Tocopherol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.09	14.62 ng	572182	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	dl-.alpha.-Tocopherol	430	C29H50O2	010191-41-0	94
2		Vitamin E	430	C29H50O2	000059-02-9	92
3		(+)-.gamma.-Tocopherol, 0-methyl-	430	C29H50O2	1000374-72-2	91
4		Propanenitrile, 3-[methyl(4-nitr...	205	C10H11N3O2	097473-81-9	47
5		1H-Purin-2-amine, 6-methoxy-	165	C6H7N5O	020535-83-5	42



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
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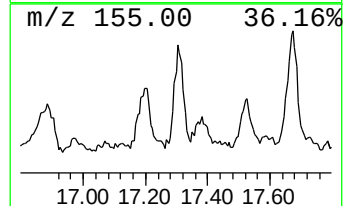
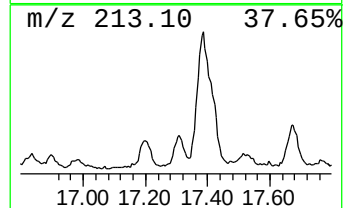
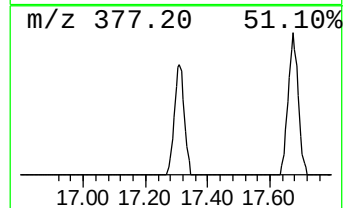
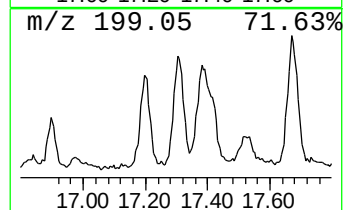
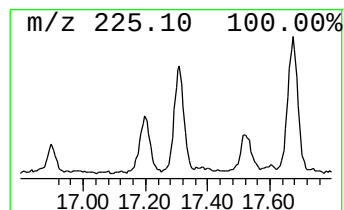
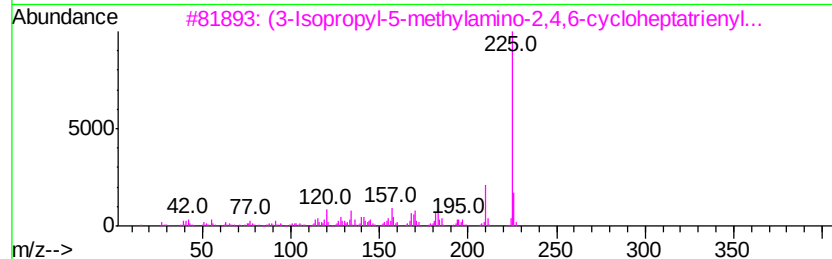
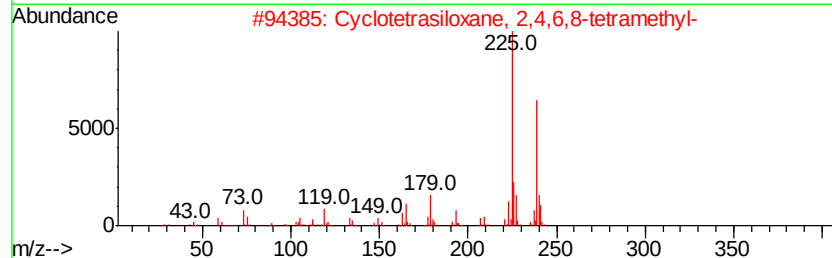
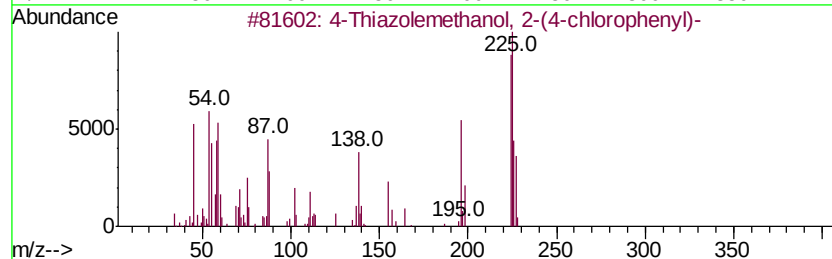
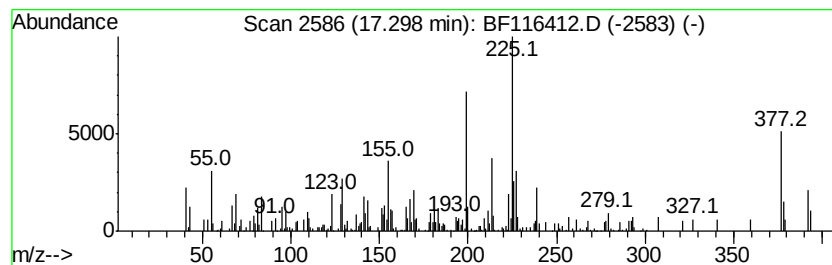
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown17.30 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.30	4.94 ng	193133	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Thiazolemethanol, 2-(4-chlorop...	225	C10H8ClNOS	036093-99-9	35
2		Cyclotetrasiloxane, 2,4,6,8-tetr...	240	C4H16O4Si4	002370-88-9	25
3		(3-Isopropyl-5-methylamino-2,4,6...	225	C14H15N3	014204-00-3	18
4		[1,2,4]Triazolo[1,5-a]pyrimidine...	225	C12H11N5	1000320-01-2	18
5		4-Chloro-1,3-dimethyl-1H-pyrazol...	225	C9H8ClN3O2	1000293-90-4	18



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
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Sample : K4223-21
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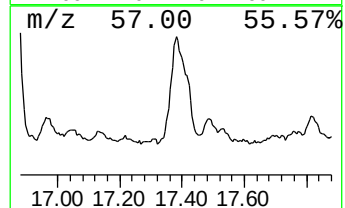
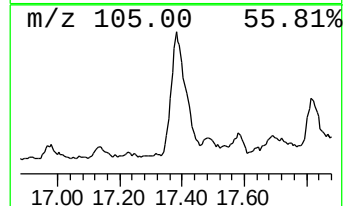
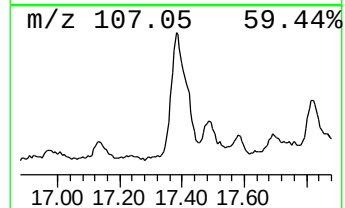
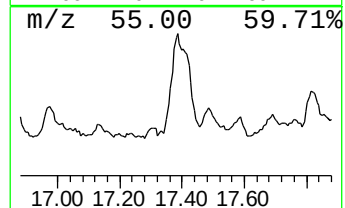
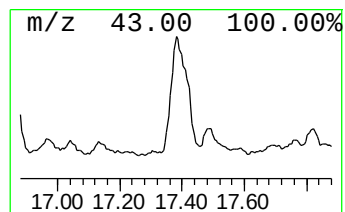
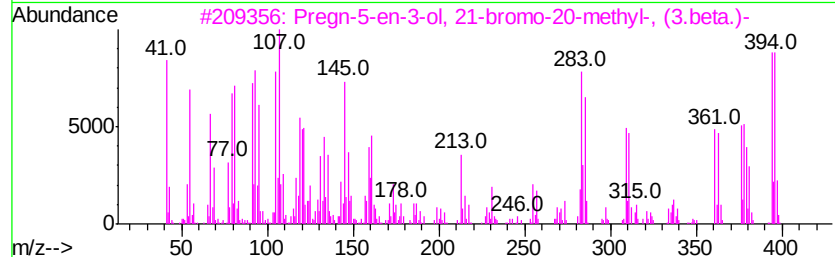
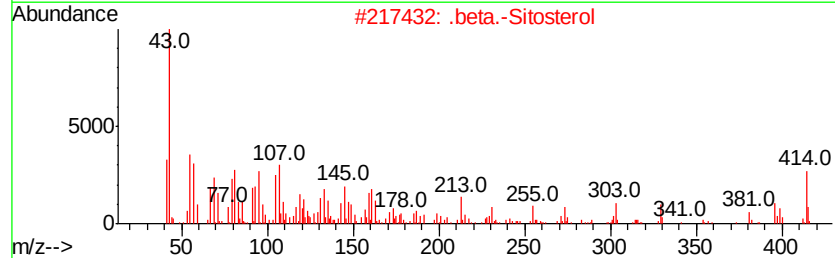
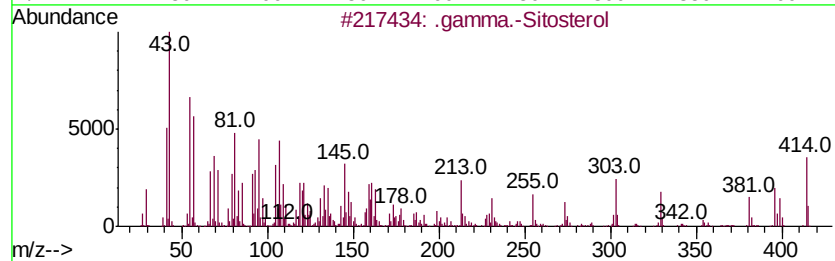
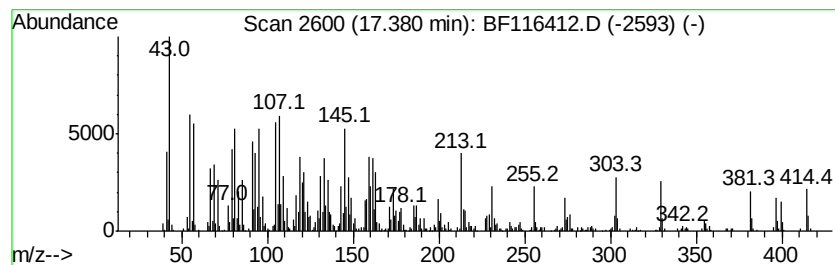
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 .gamma.-Sitosterol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.38	72.39 ng	2832620	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	99
3		Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	64
4		Ergost-5-en-3-ol, (3.beta.)-	400	C28H48O	004651-51-8	52
5		Isocholesteryl methyl ether	400	C28H48O	029944-53-4	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
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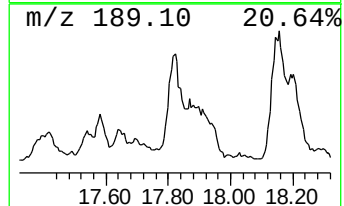
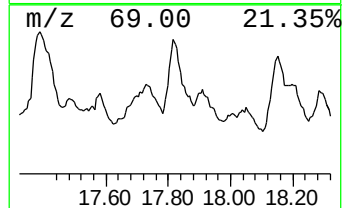
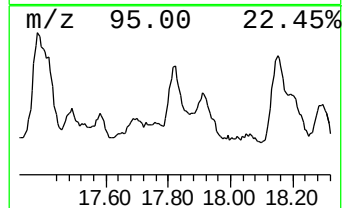
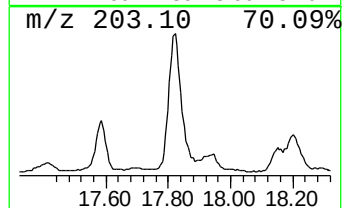
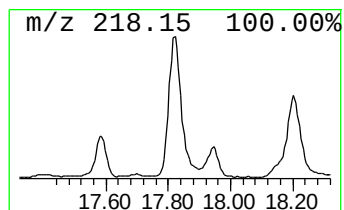
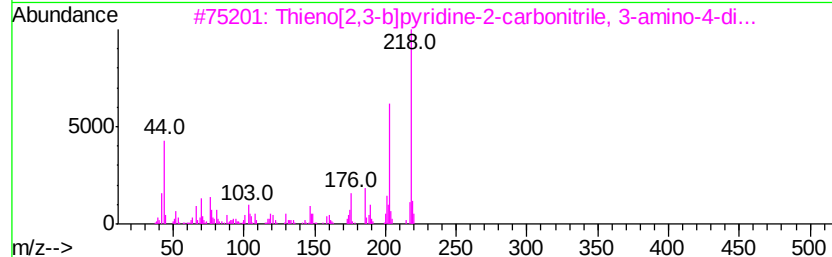
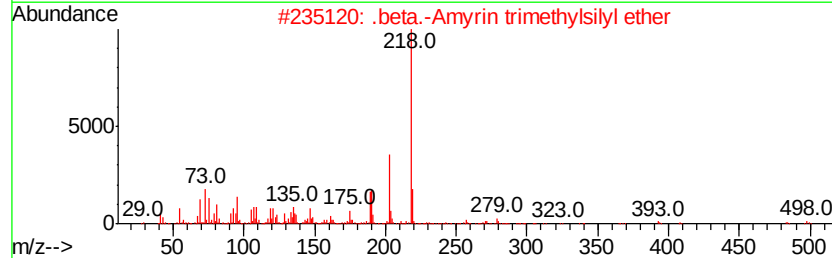
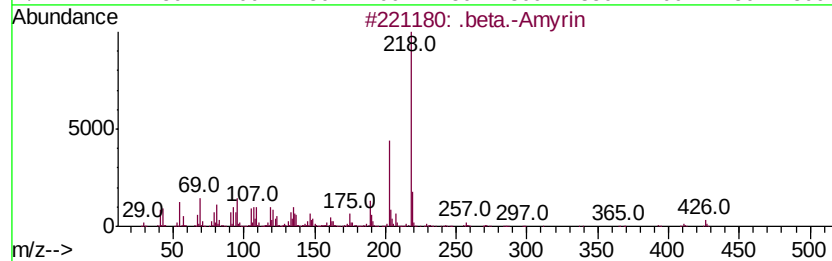
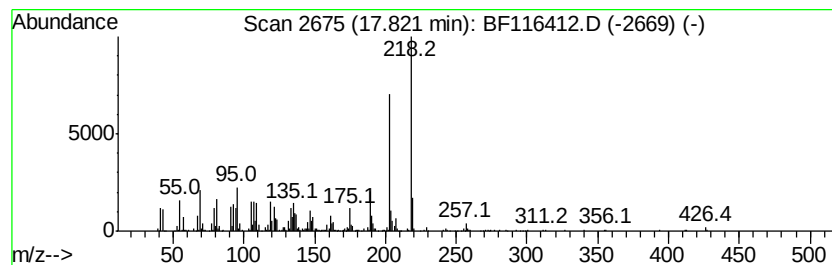
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 .beta.-Amyrin Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.82	23.47 ng	918353	Perylene-d12	15.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		.beta.-Amyrin	426 C30H50O	000559-70-6 98
2		.beta.-Amyrin trimethylsilyl ether	498 C33H58OSi	001721-67-1 80
3		Thieno[2,3-b]pyridine-2-carbonit...	218 C10H10N4S	147992-88-9 64
4		2(1H)Naphthalenone, 3,5,6,7,8,8a...	218 C15H22O	1000188-66-5 64
5		6,7,8,9-Tetrahydrodibenzo[b,d]fu...	260 C15H16O4	1000196-34-3 62



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Data File : BF116412.D
Acq On : 20 Aug 2019 2:26
Operator : HP/JU
Sample : K4223-21
Misc :
ALS Vial : 21 Sample Multiplier: 1

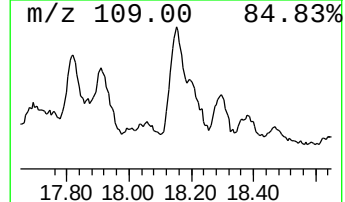
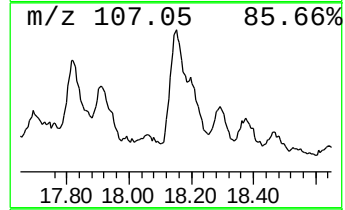
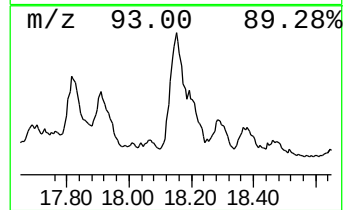
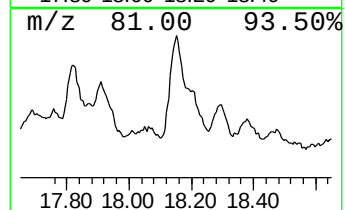
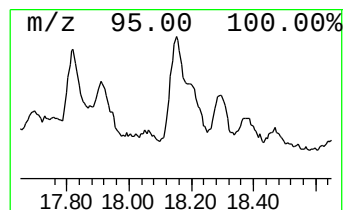
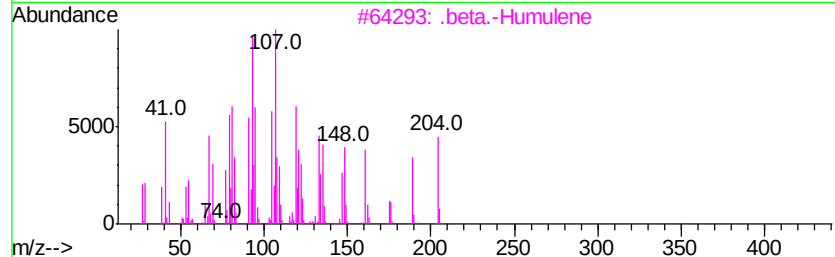
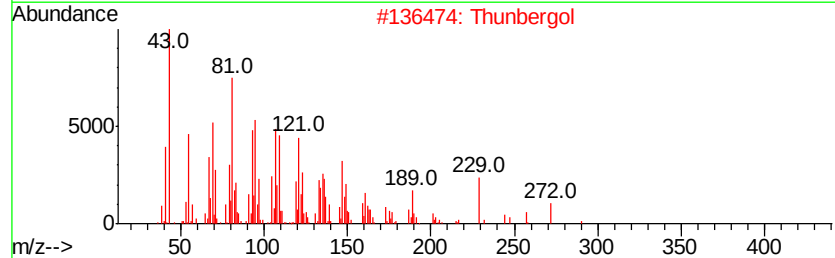
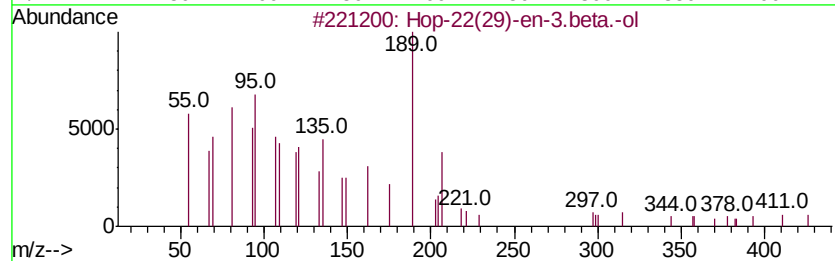
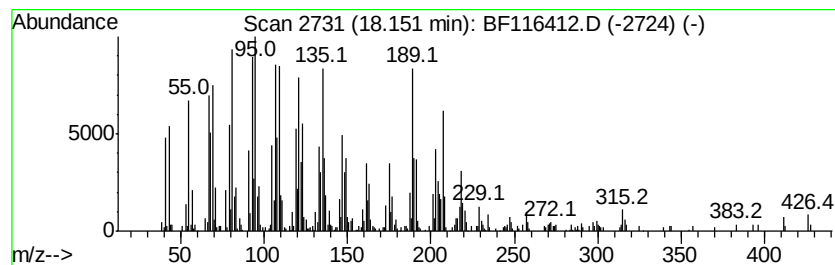
Instrument :
BNA_F
ClientSampled :
SP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Hop-22(29)-en-3.beta.-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.15	30.38 ng	1188620	Perylene-d12	15.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Hop-22(29)-en-3.beta.-ol	426 C30H50O	058801-23-3 92
2		Thunbergol	290 C20H34O	025269-17-4 90
3		.beta.-Humulene	204 C15H24	000116-04-1 78
4		Longifolenaldehyde	220 C15H24O	019890-84-7 64
5		Guaia-9,11-diene	204 C15H24	1000374-19-8 52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081919\
Data File : BF116412.D
Acq On : 20 Aug 2019 2:26
Operator : HP/JU
Sample : K4223-21
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-11

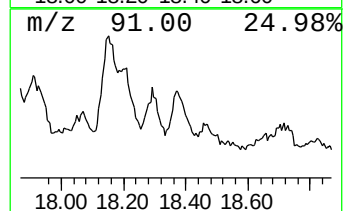
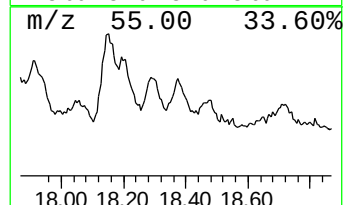
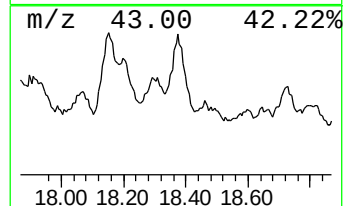
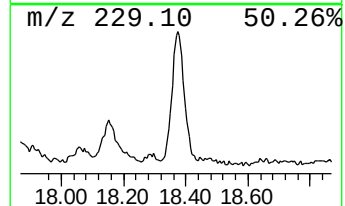
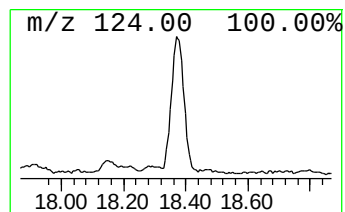
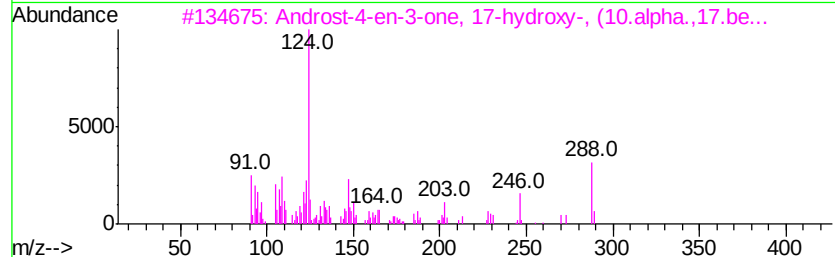
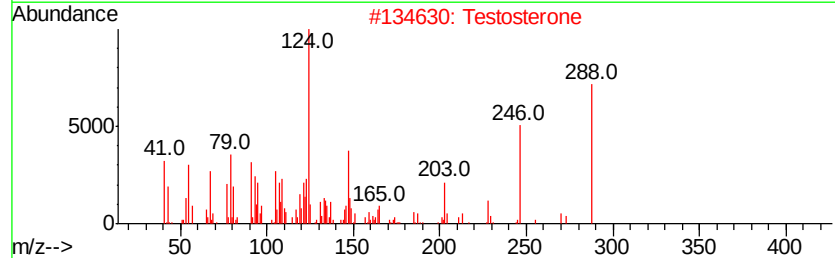
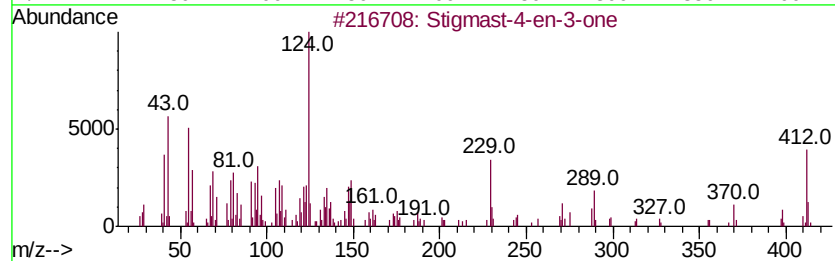
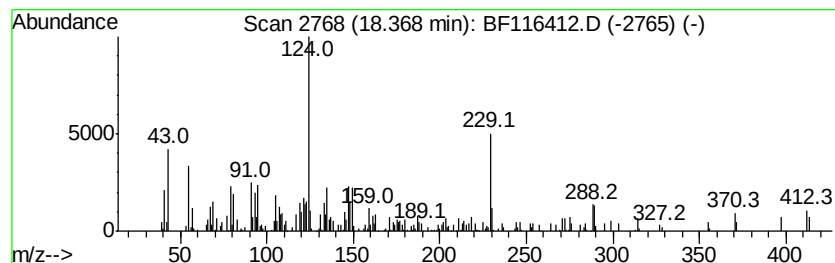
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Stigmast-4-en-3-one Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	11.93 ng	466678	Perylene-d12	15.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Stigmast-4-en-3-one	412	C29H48O	001058-61-3	89
2		Testosterone	288	C19H28O2	000058-22-0	58
3		Androst-4-en-3-one, 17-hydroxy-,...	288	C19H28O2	000604-39-7	41
4		Cholest-5-en-3-one, 4,4-dimethyl-	412	C29H48O	002220-42-0	38
5		Orcinol	124	C7H8O2	000504-15-4	30



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 Operator : HP/JU
 Sample : K4223-21
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-11

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073019.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
unknown6.58	6.58	68.7	ng	2080720	1	6.80	605466	20.0
Naphthalene, 1,2,...	9.95	3.7	ng	180831	3	9.83	988406	20.0
n-Hexadecanoic acid	11.85	20.6	ng	1124800	4	11.32	1094760	20.0
17-Norkaur-15-ene...	12.00	4.6	ng	253243	4	11.32	1094760	20.0
Phytol	12.45	7.1	ng	390216	4	11.32	1094760	20.0
Octadecanoic acid	12.63	4.2	ng	228246	4	11.32	1094760	20.0
2-Phenanthrenol, ...	13.33	5.2	ng	199242	5	13.96	768594	20.0
1-Dodecanol, 2-oc...	13.79	12.6	ng	482417	5	13.96	768594	20.0
Sulfurous acid, d...	14.41	19.8	ng	760273	5	13.96	768594	20.0
Squalene	14.78	6.1	ng	237142	6	15.42	782568	20.0
10-Heneicosene (c,t)	15.06	5.1	ng	198349	6	15.42	782568	20.0
Eicosane	15.82	39.7	ng	1553260	6	15.42	782568	20.0
n-Tetracosanol-1	15.87	6.8	ng	267883	6	15.42	782568	20.0
dl-.alpha.-Tocoph...	16.09	14.6	ng	572182	6	15.42	782568	20.0
unknown17.30	17.30	4.9	ng	193133	6	15.42	782568	20.0
.gamma.-Sitosterol	17.38	72.4	ng	2832620	6	15.42	782568	20.0
.beta.-Amyrin	17.82	23.5	ng	918353	6	15.42	782568	20.0
Hop-22(29)-en-3.b...	18.15	30.4	ng	1188620	6	15.42	782568	20.0
Stigmast-4-en-3-one	18.37	11.9	ng	466678	6	15.42	782568	20.0