

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
 Data File : BF103232.D
 Acq On : 26 Feb 2018 21:57
 Operator : SJ/JU
 Sample : J1691-53 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B14

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

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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	5.093	508	511	519	rBV	33153	45479	2.56%	0.185%
2	5.492	575	579	592	rBV	1767626	1714384	96.48%	6.977%
3	6.504	747	751	771	rBV	1791131	1776931	100.00%	7.231%
4	6.645	771	775	778	rVV	1929408	1708494	96.15%	6.953%
5	6.698	781	784	789	rVB3	19548	20024	1.13%	0.081%
6	6.875	810	814	821	rBV	1157086	981968	55.26%	3.996%
7	7.028	836	840	848	rVB	1513661	1300848	73.21%	5.294%
8	7.434	905	909	919	rBV	1086376	1033956	58.19%	4.208%
9	8.157	1028	1032	1035	rBV	1411446	1234468	69.47%	5.024%
10	8.563	1098	1101	1104	rBV	24615	20310	1.14%	0.083%
11	8.710	1124	1126	1128	rBV	72352	57885	3.26%	0.236%
12	8.733	1128	1130	1134	rVB	47858	38898	2.19%	0.158%
13	8.869	1150	1153	1156	rBV	47779	45988	2.59%	0.187%
14	8.969	1167	1170	1173	rBV	40373	34120	1.92%	0.139%
15	9.163	1201	1203	1210	rVB	28821	24783	1.39%	0.101%
16	9.228	1210	1214	1224	rBV	1939748	1628535	91.65%	6.627%
17	9.304	1224	1227	1229	rVV	63732	61821	3.48%	0.252%
18	9.333	1229	1232	1236	rVB3	26433	29435	1.66%	0.120%
19	9.498	1256	1260	1265	rBV2	37984	59468	3.35%	0.242%
20	9.569	1268	1272	1275	rBV	49924	60468	3.40%	0.246%
21	9.622	1278	1281	1283	rVB	109817	93765	5.28%	0.382%
22	9.686	1288	1292	1294	rBV3	26774	29924	1.68%	0.122%
23	9.775	1303	1307	1312	rBV	93816	91261	5.14%	0.371%
24	9.828	1312	1316	1320	rVB	86483	88755	4.99%	0.361%
25	9.916	1327	1331	1334	rBV	1274951	1147740	64.59%	4.671%
26	9.945	1334	1336	1340	rVB	212075	165212	9.30%	0.672%
27	10.016	1344	1348	1353	rBV4	16500	23364	1.31%	0.095%
28	10.069	1353	1357	1359	rBV3	27307	31230	1.76%	0.127%
29	10.116	1362	1365	1369	rVB2	78131	79922	4.50%	0.325%
30	10.151	1369	1371	1376	rVB2	42109	42949	2.42%	0.175%
31	10.227	1380	1384	1387	rBV3	31586	36476	2.05%	0.148%
32	10.257	1387	1389	1395	rVB2	21050	24645	1.39%	0.100%
33	10.327	1395	1401	1405	rBV2	111416	129467	7.29%	0.527%
34	10.463	1421	1424	1430	rVB	166903	161117	9.07%	0.656%

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Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
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35	10.551	1433	1439	1441	rVV4	62839	78018	4.39%	0.317%
36	10.622	1448	1451	1456	rVB	321224	270483	15.22%	1.101%
37	10.704	1462	1465	1471	rBV	895196	813543	45.78%	3.311%
38	10.804	1479	1482	1492	rVB	125698	173471	9.76%	0.706%
39	10.892	1495	1497	1500	rBV4	26994	27385	1.54%	0.111%
40	11.010	1511	1517	1520	rBV2	46400	74653	4.20%	0.304%
41	11.092	1528	1531	1533	rVB3	24254	20243	1.14%	0.082%
42	11.133	1534	1538	1540	rBV4	25305	35139	1.98%	0.143%
43	11.210	1548	1551	1555	rVB4	16745	19076	1.07%	0.078%
44	11.251	1555	1558	1561	rBV2	107390	110537	6.22%	0.450%
45	11.280	1561	1563	1565	rVV	60960	61994	3.49%	0.252%
46	11.298	1565	1566	1570	rVB	63945	52278	2.94%	0.213%
47	11.404	1580	1584	1586	rBV	1090926	976757	54.97%	3.975%
48	11.427	1586	1588	1593	rVV	927676	780202	43.91%	3.175%
49	11.480	1594	1597	1601	rVV	241856	223203	12.56%	0.908%
50	11.516	1601	1603	1608	rVB2	37443	43360	2.44%	0.176%
51	11.639	1620	1624	1628	rBV	72589	93030	5.24%	0.379%
52	11.680	1628	1631	1636	rVB2	69053	69913	3.93%	0.285%
53	11.904	1666	1669	1672	rBV	99357	118127	6.65%	0.481%
54	11.933	1672	1674	1678	rVB	135815	124012	6.98%	0.505%
55	11.980	1680	1682	1685	rBV	53144	59251	3.33%	0.241%
56	12.021	1685	1689	1691	rVV	194959	213771	12.03%	0.870%
57	12.039	1691	1692	1697	rVB	137133	106445	5.99%	0.433%
58	12.086	1697	1700	1703	rBV	85713	72180	4.06%	0.294%
59	12.145	1707	1710	1713	rBV3	67424	73292	4.12%	0.298%
60	12.198	1717	1719	1722	rBV	65312	54333	3.06%	0.221%
61	12.233	1723	1725	1726	rBV2	30895	22660	1.28%	0.092%
62	12.257	1726	1729	1731	rVB	114437	107816	6.07%	0.439%
63	12.457	1760	1763	1765	rBV	74738	83325	4.69%	0.339%
64	12.474	1765	1766	1769	rVB	44324	37526	2.11%	0.153%
65	12.621	1787	1791	1794	rBV	1127221	981715	55.25%	3.995%
66	12.651	1794	1796	1800	rVB2	74267	84148	4.74%	0.342%
67	12.851	1824	1830	1836	rBV	918233	1136801	63.98%	4.626%
68	12.986	1850	1853	1856	rVB	1240362	1017791	57.28%	4.142%
69	13.180	1883	1886	1888	rVB	167105	146424	8.24%	0.596%
70	13.245	1894	1897	1899	rBV	173126	147005	8.27%	0.598%
71	13.874	2002	2004	2007	rVB2	183325	152684	8.59%	0.621%

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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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Peak separation: 5

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72	14.057	2030	2035	2037	rBV2	809496	1106531	62.27%	4.503%
73	14.080	2037	2039	2045	rVB	388492	341979	19.25%	1.392%
74	15.557	2287	2290	2295	rVB	335382	437909	24.64%	1.782%

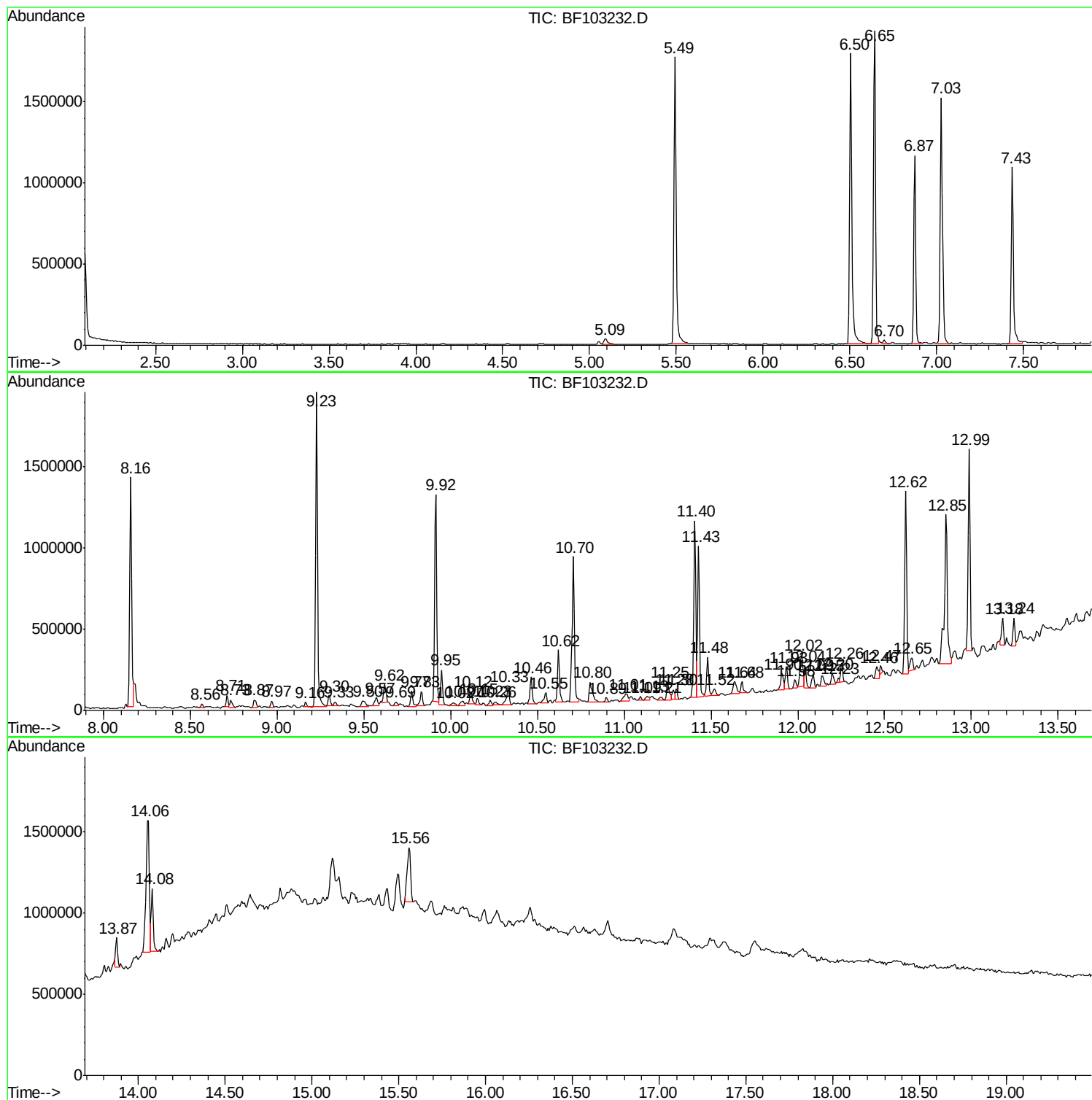
Sum of corrected areas: 24573100

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Instrument :
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.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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Sample : J1691-53 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

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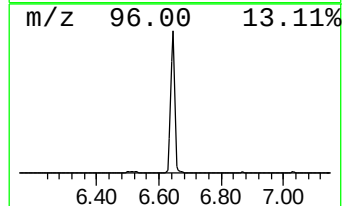
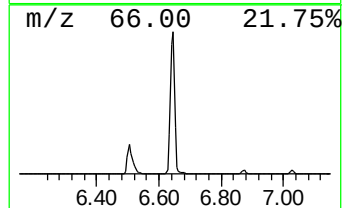
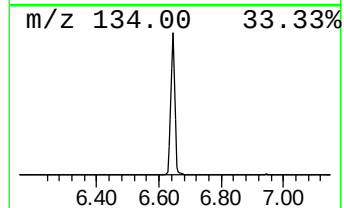
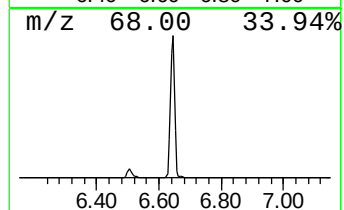
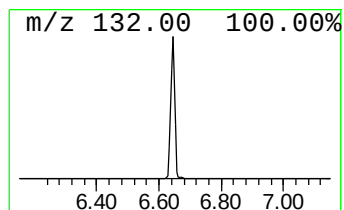
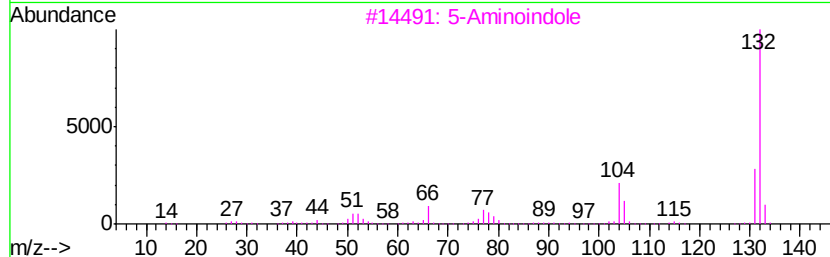
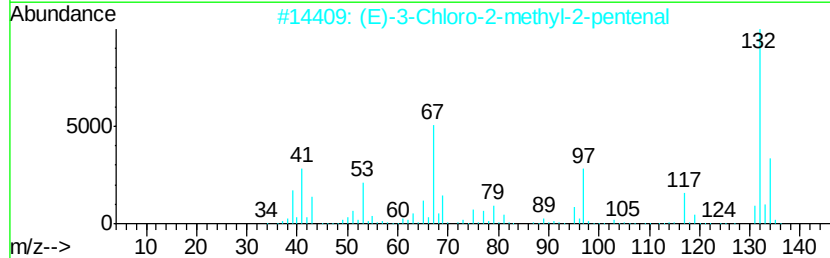
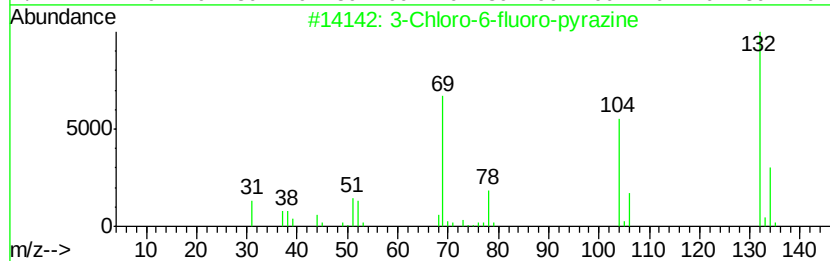
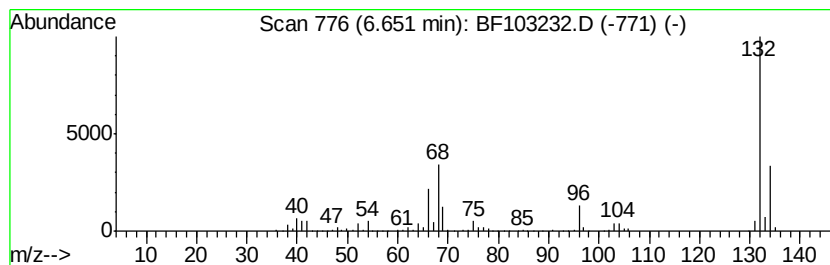
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	34.80 ng	1708490	1,4-Dichlorobenzene-d4	6.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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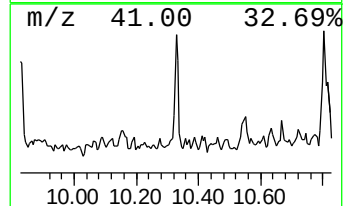
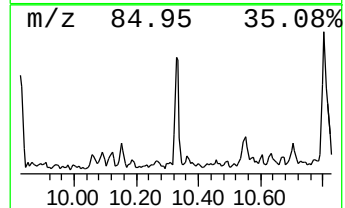
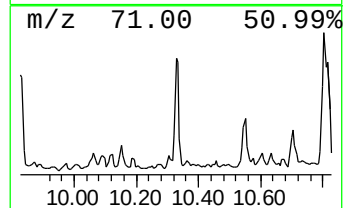
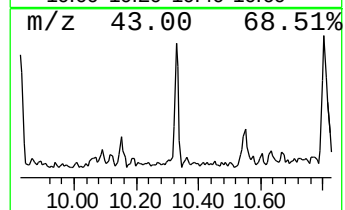
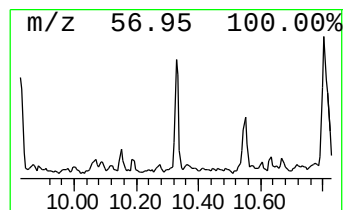
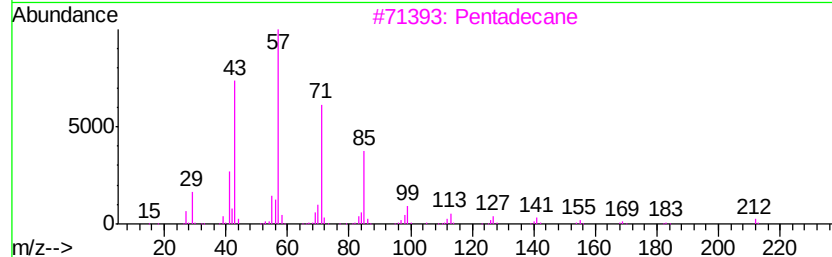
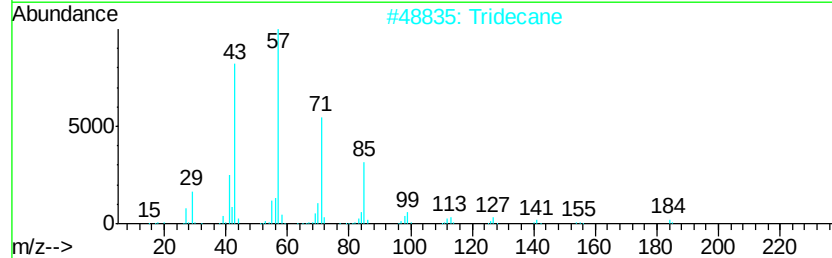
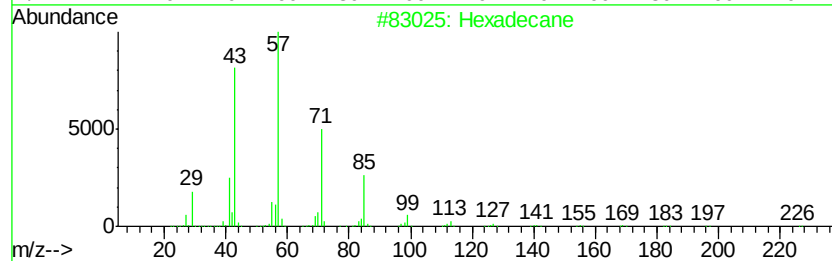
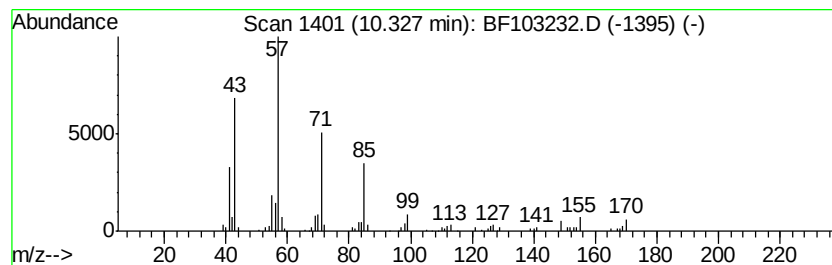
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	2.26 ng	129467	Acenaphthene-d10	9.92
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	83
2	Tridecane	184 C13H28	000629-50-5	72
3	Pentadecane	212 C15H32	000629-62-9	64
4	Sulfurous acid, butyl nonyl ester	264 C13H28O3S	1000309-17-6	64
5	Tetradecane	198 C14H30	000629-59-4	58



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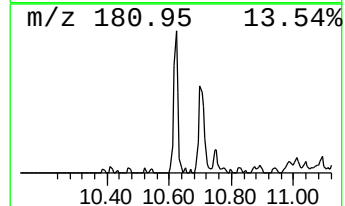
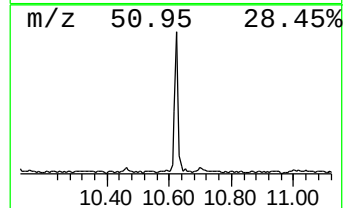
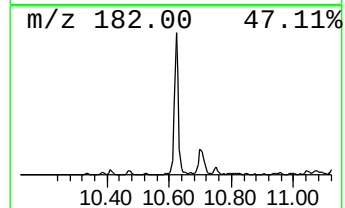
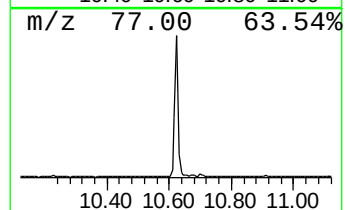
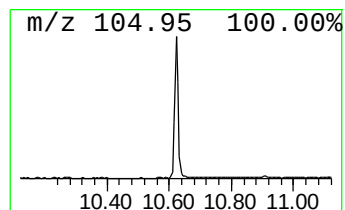
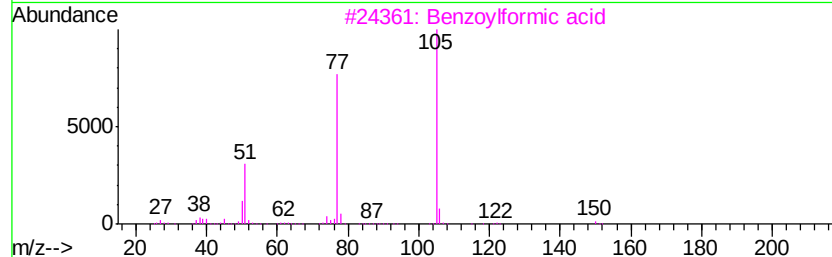
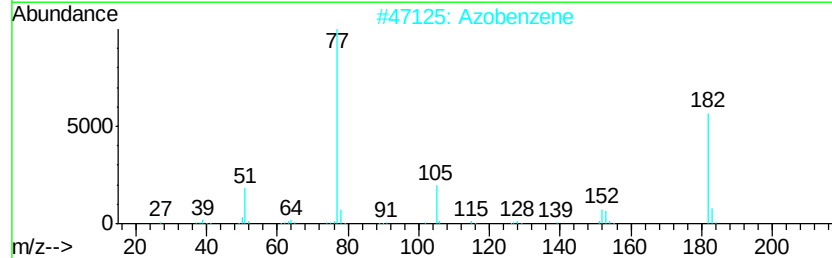
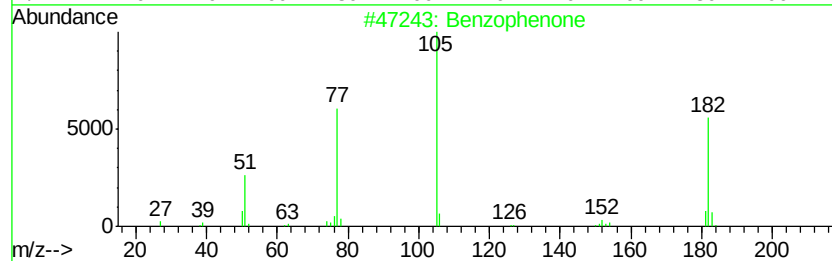
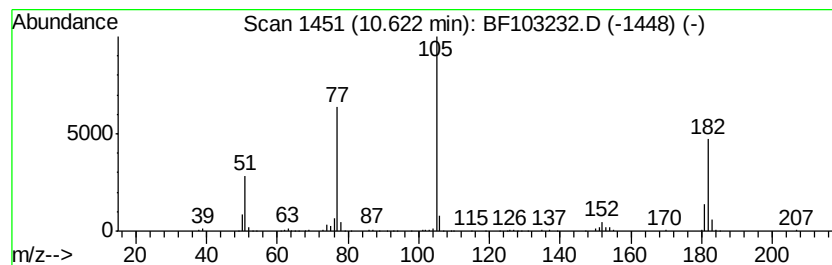
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.62	4.71 ng	270483	Acenaphthene-d10		9.92	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzophenone	182	C13H10O	000119-61-9	96
2		Azobenzene	182	C12H10N2	000103-33-3	58
3		Benzoylformic acid	150	C8H6O3	000611-73-4	49
4		Phenyl 4-pyridyl ketone	183	C12H9NO	014548-46-0	47
5		2-(2-Oxo-2-phenyl-ethyl)-malonon...	184	C11H8N2O	1000296-76-9	47



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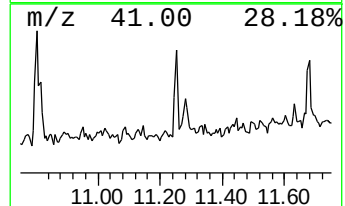
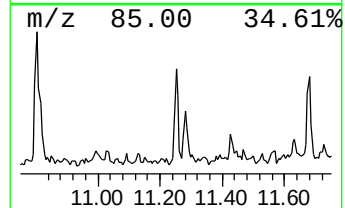
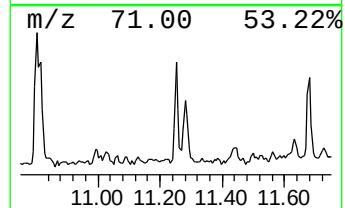
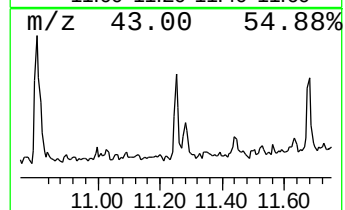
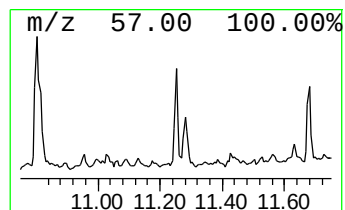
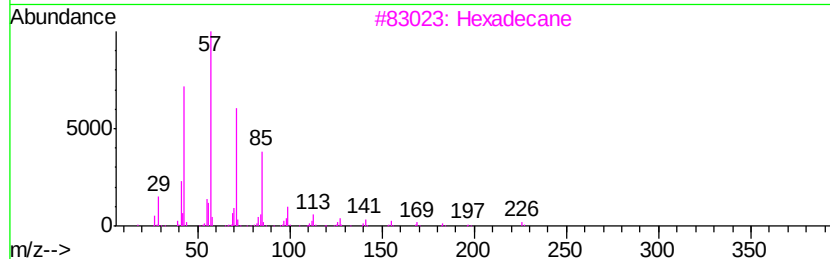
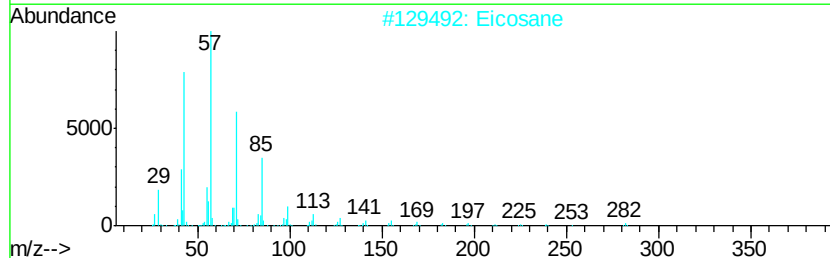
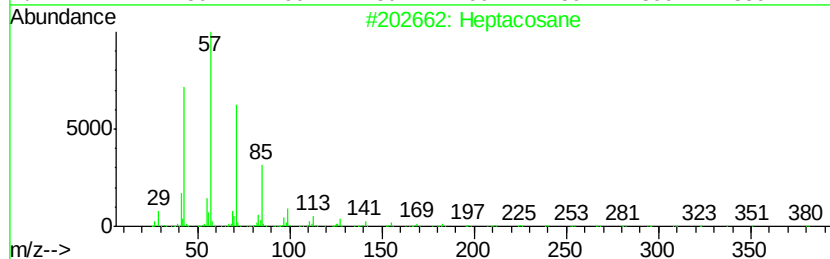
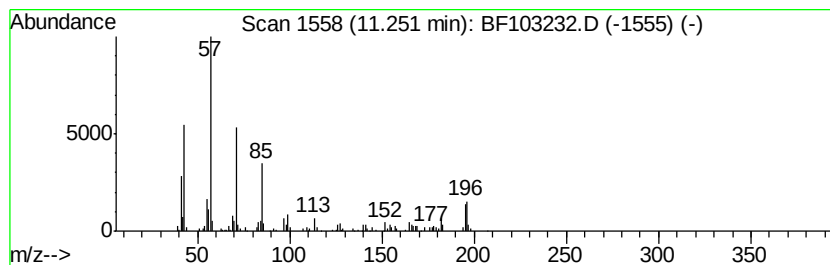
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.25	2.26 ng	110537	Phenanthrene-d10	11.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptacosane	380 C27H56	000593-49-7	52
2	Eicosane	282 C20H42	000112-95-8	52
3	Hexadecane	226 C16H34	000544-76-3	52
4	Nonane, 3-methyl-	142 C10H22	005911-04-6	46
5	Dodecane, 1-iodo-	296 C12H25I	004292-19-7	46



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Sample : J1691-53 2X
Misc :
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Instrument :
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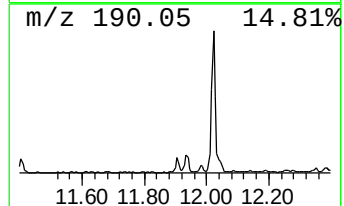
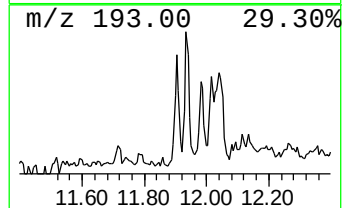
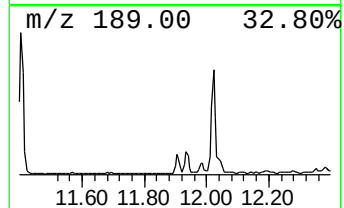
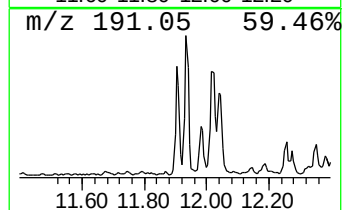
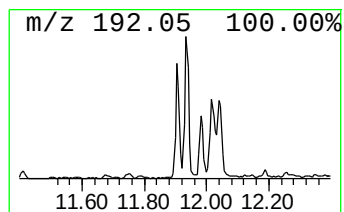
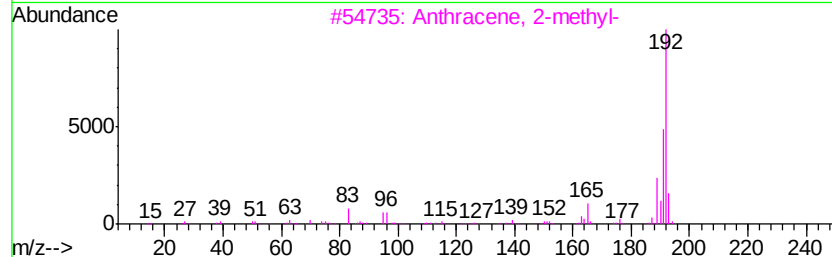
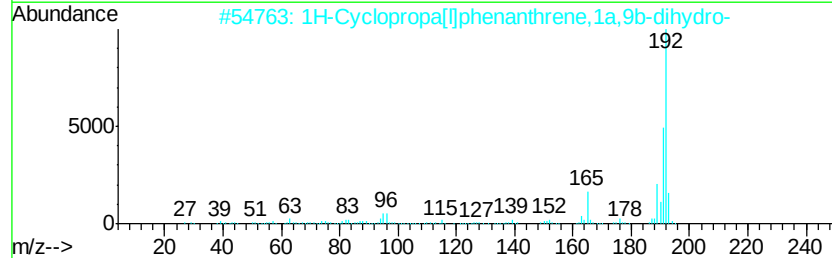
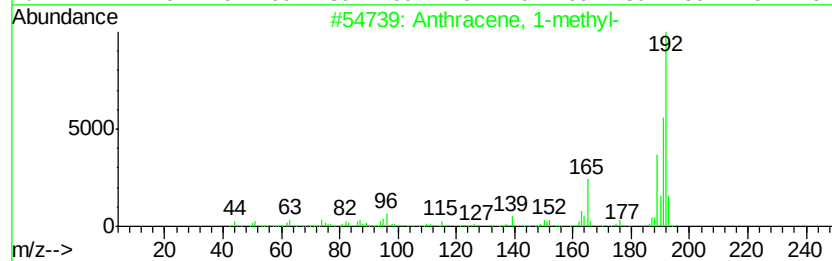
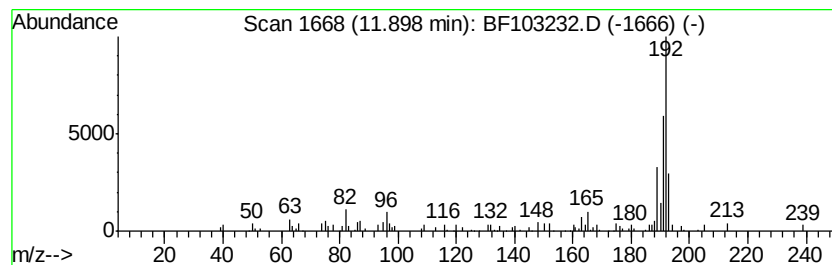
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Anthracene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	2.42 ng	118127	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91



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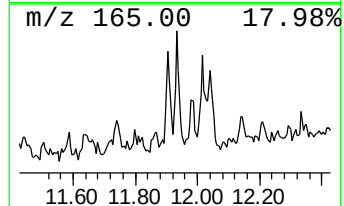
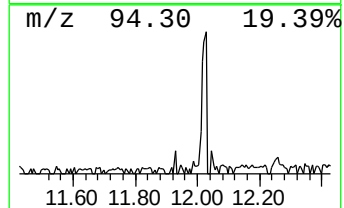
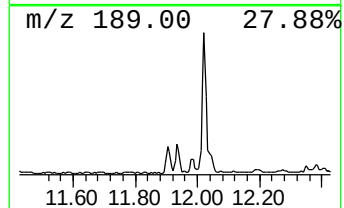
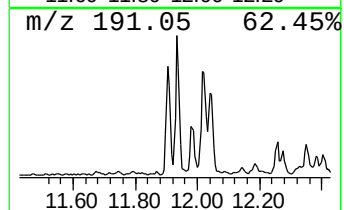
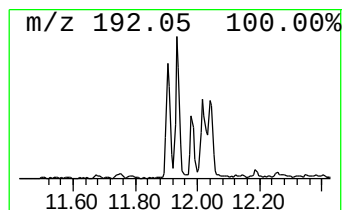
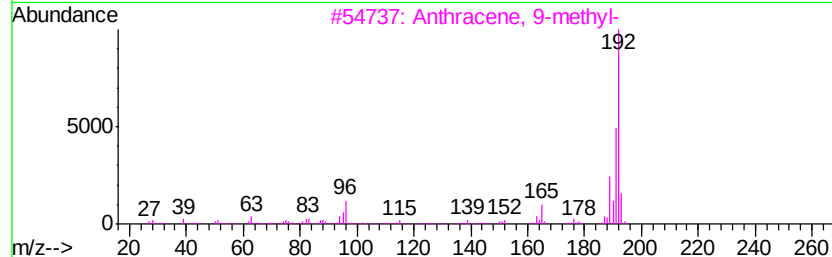
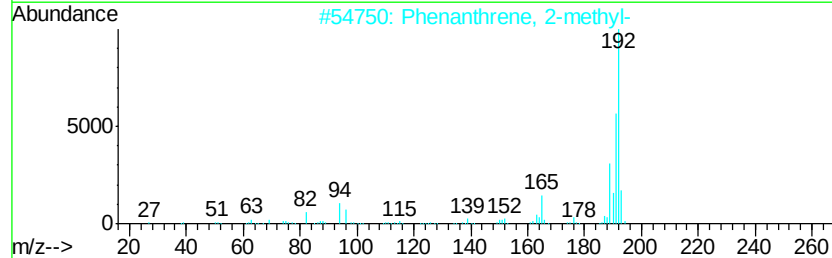
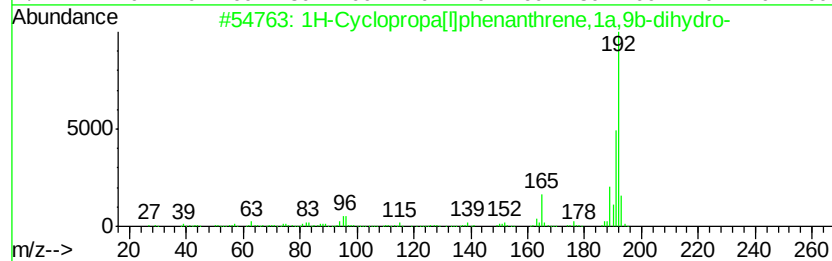
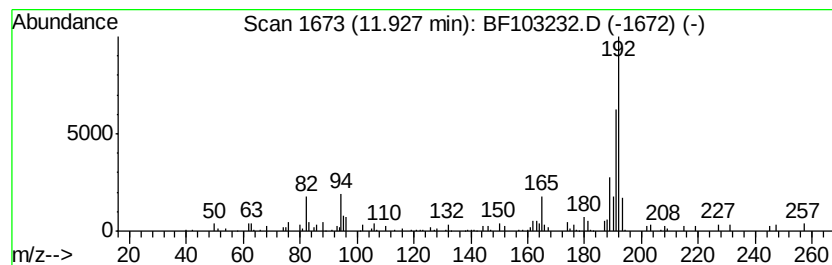
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Cyclopropa[l]phenanthren... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	2.54 ng	124012	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	96
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



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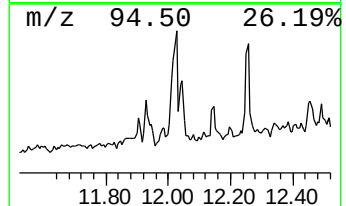
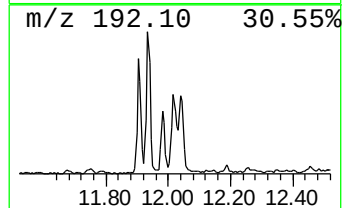
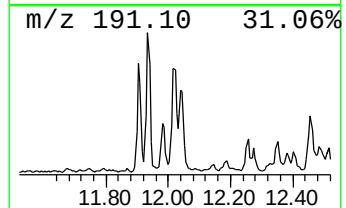
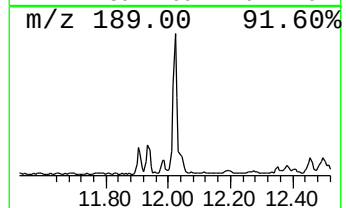
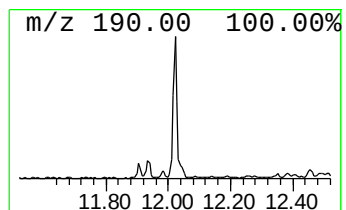
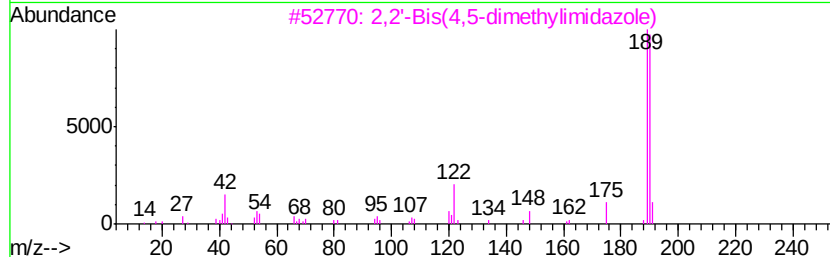
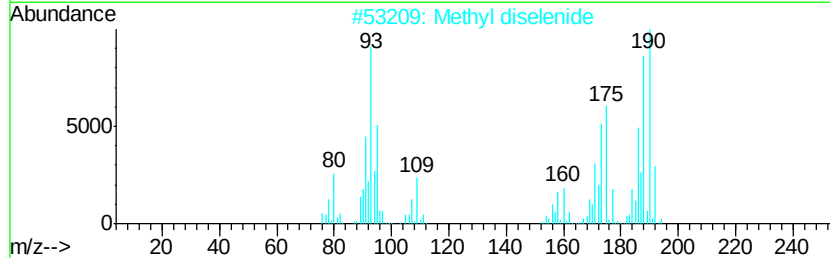
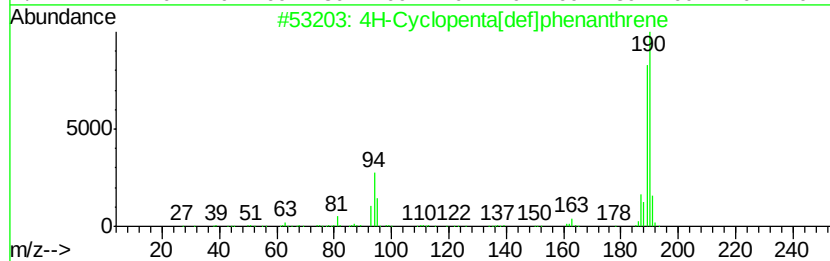
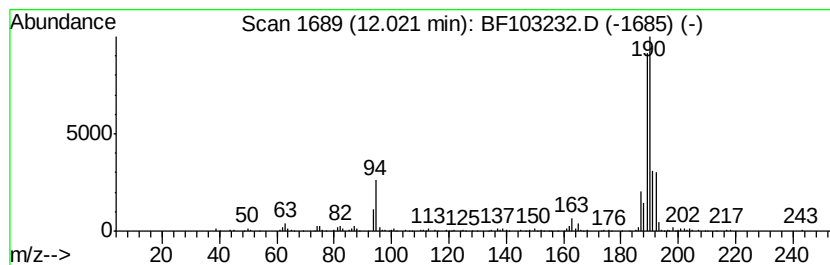
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.02	4.38 ng	213771	Phenanthrene-d10	11.40		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2		Methyl diselenide	190	C2H6Se2	007101-31-7	49
3		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
4		1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25
5		Pyrimidine, 2-chloro-5-phenyl-	190	C10H7ClN2	022536-62-5	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
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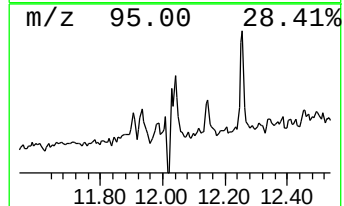
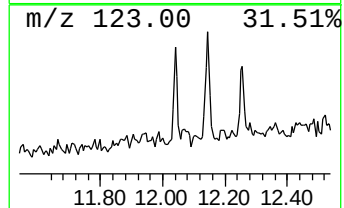
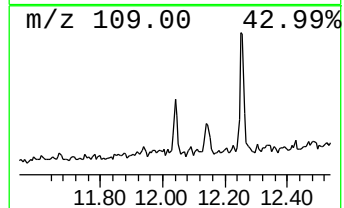
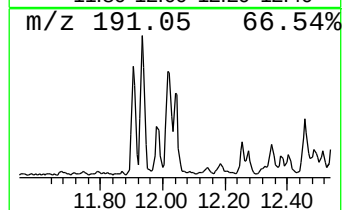
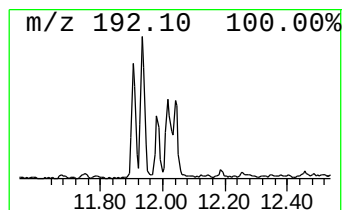
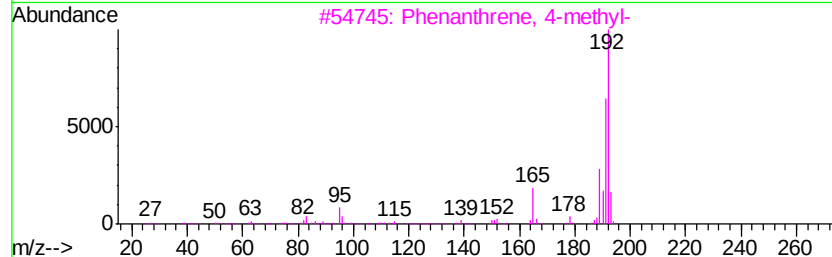
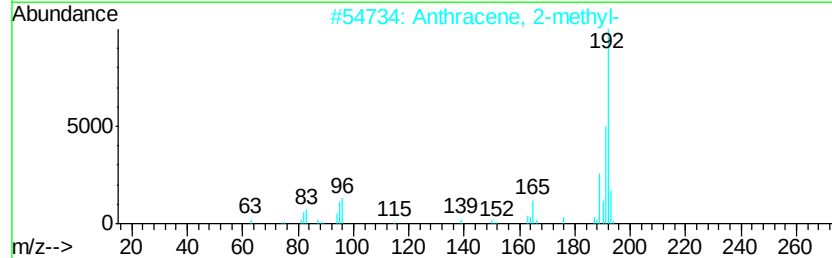
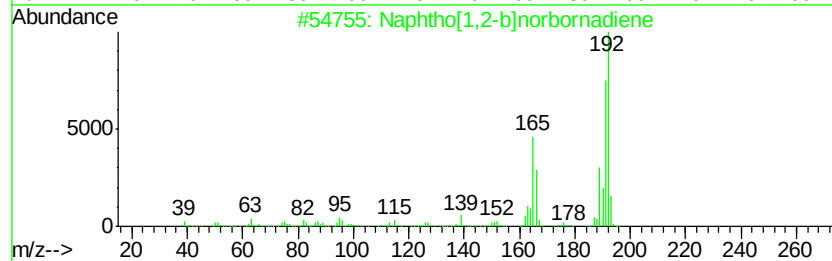
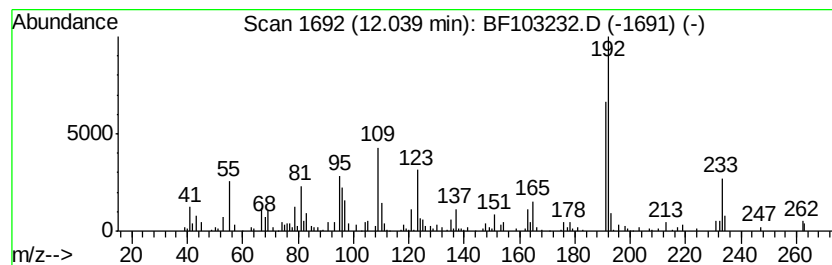
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown12.04 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	2.18 ng	106445	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	46
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	43
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	43
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	43
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	43



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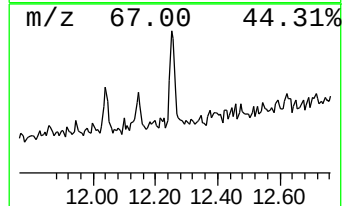
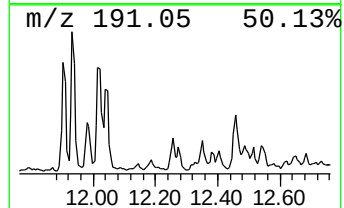
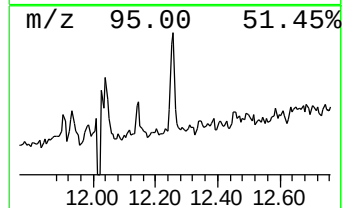
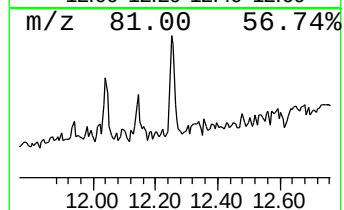
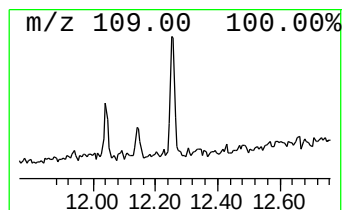
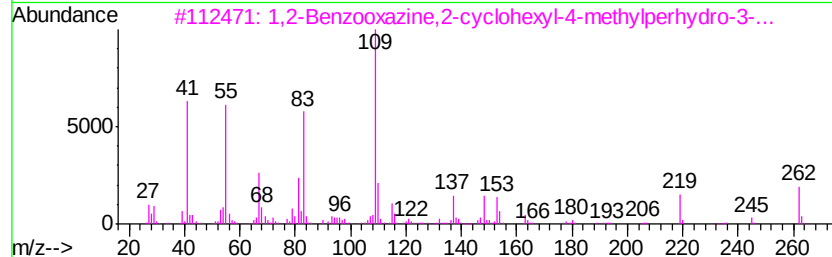
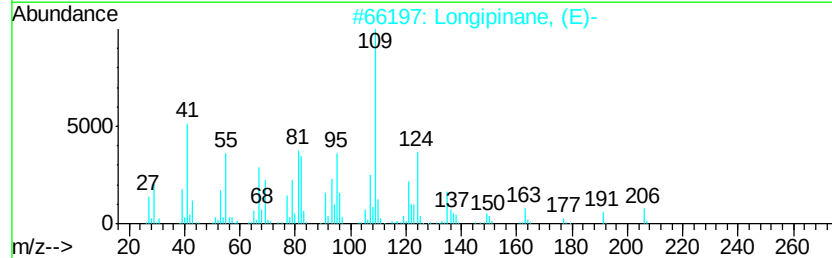
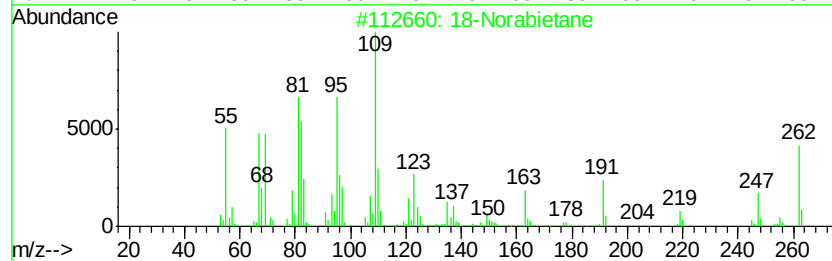
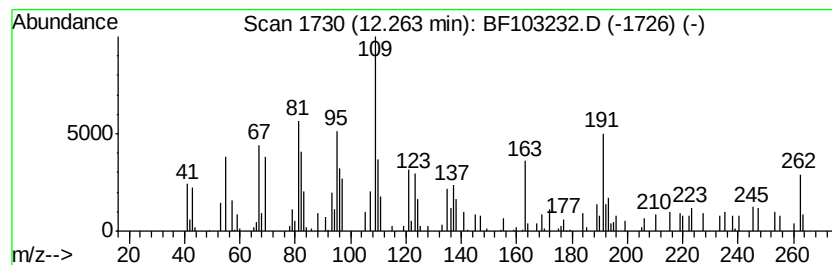
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TIC Integration Parameters: LSCINT.P

Peak Number 11 18-Norabietane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.26	2.21 ng	107816	Phenanthrene-d10	11.40	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	18-Norabietane	262	C19H34	1000293-16-6	93
2	Longipinane, (E)-	206	C15H26	1000156-14-5	43
3	1,2-Benzooxazine,2-cyclohexyl-4-...	262	C16H26N2O	037898-39-8	30
4	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	22
5	1,2-Dioctylcyclopropene	264	C19H36	001089-40-3	22



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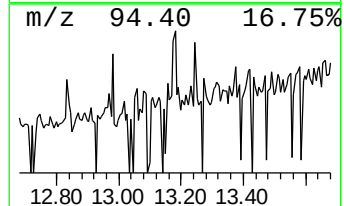
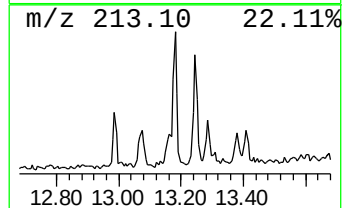
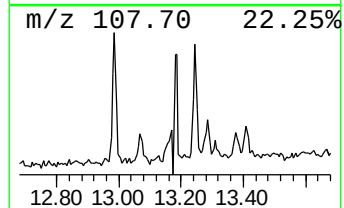
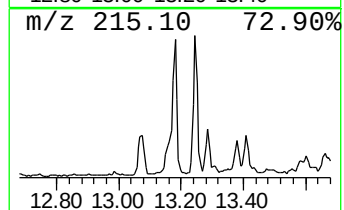
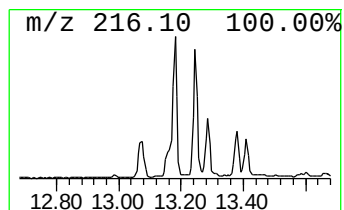
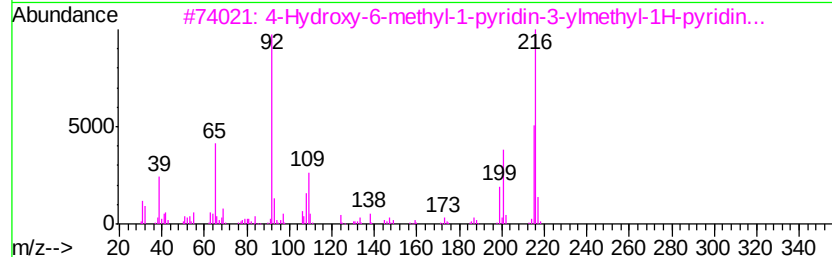
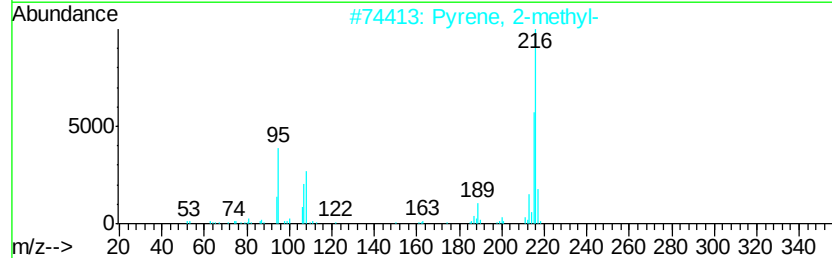
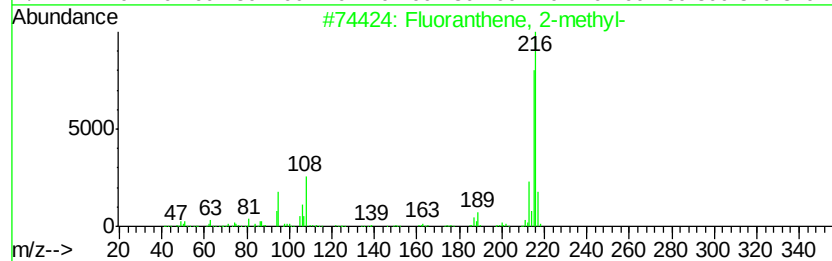
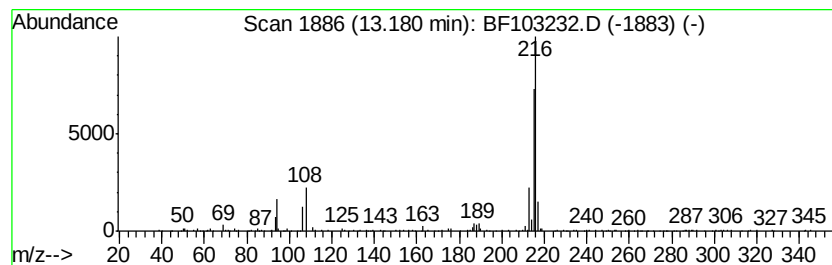
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Fluoranthene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	2.65 ng	146424	Chrysene-d12	14.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
2		Pyrene, 2-methyl-	216	C17H12	003442-78-2	87
3		4-Hydroxy-6-methyl-1-pyridin-3-ylmethyl-1H-pyridin...	216	C12H12N2O2	1000318-25-3	74
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72
5		Pyrene, 1-methyl-	216	C17H12	002381-21-7	72



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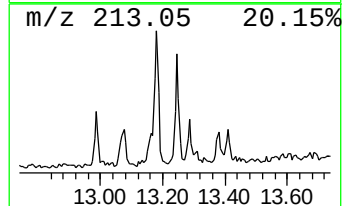
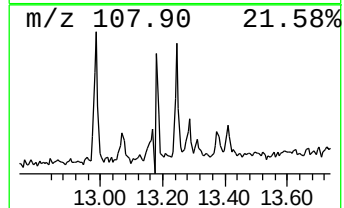
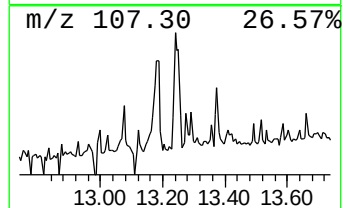
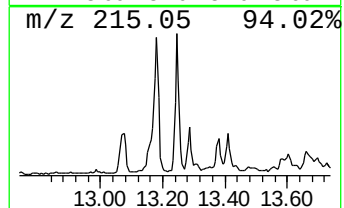
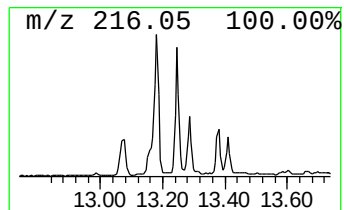
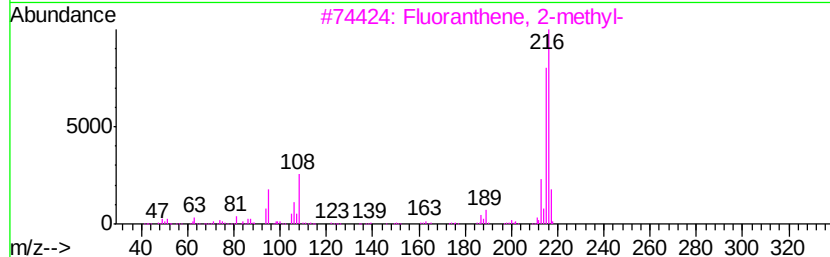
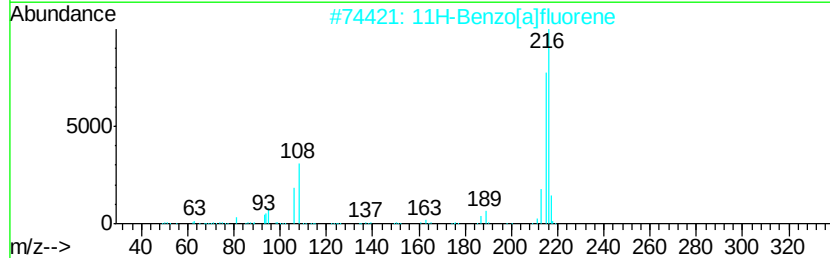
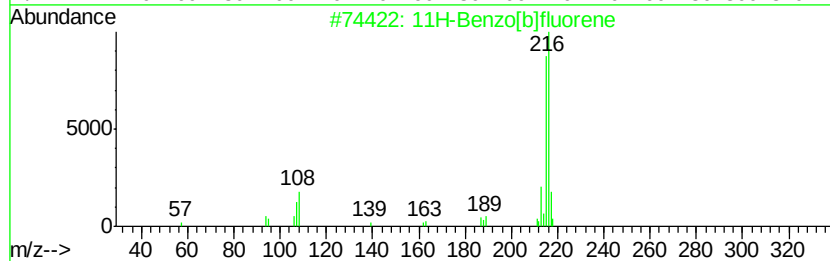
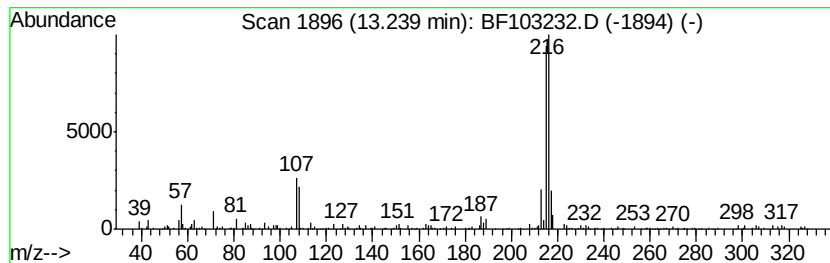
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 11H-Benzo[b]fluorene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.24	2.66 ng	147005	Chrysene-d12	14.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 93
2		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 86
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 83
4		7H-Benzo[c]fluorene	216 C17H12	000205-12-9 80
5		Pyrene, 1-methyl-	216 C17H12	002381-21-7 72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF022618\
Data File : BF103232.D
Acq On : 26 Feb 2018 21:57
Operator : SJ/JU
Sample : J1691-53 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

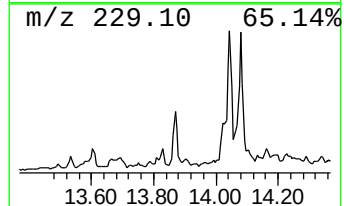
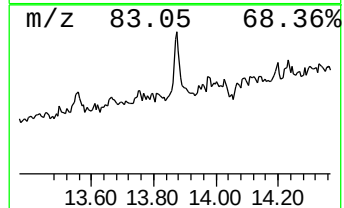
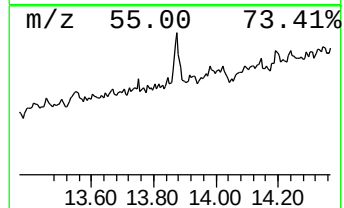
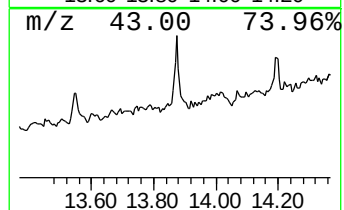
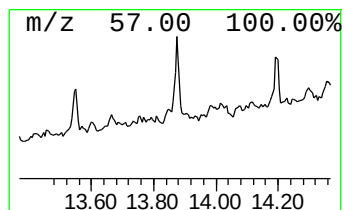
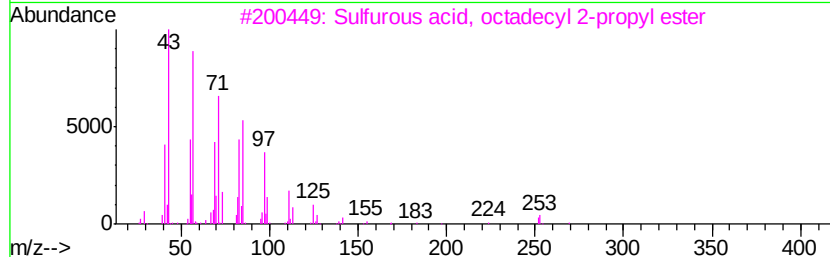
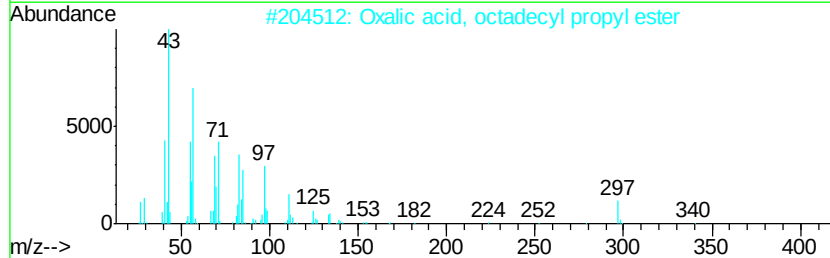
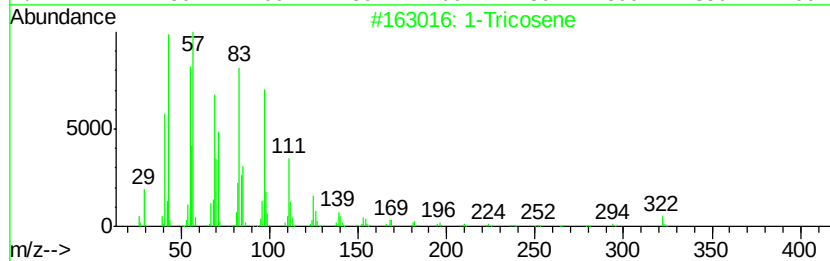
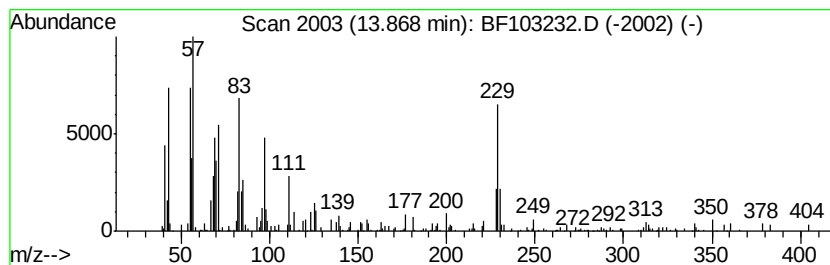
Instrument :
BNA_F
ClientSampleId :
B14

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Tricosene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	2.76 ng	152684	Chrysene-d12	14.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Tricosene	322 C23H46	018835-32-0	83
2	Oxalic acid, octadecyl propyl ester	384 C23H44O4	1000309-27-0	68
3	Sulfurous acid, octadecyl 2-prop...	376 C21H44O3S	1000309-12-7	64
4	Oxalic acid, isobutyl tetradecyl...	342 C20H38O4	1000309-37-9	64
5	Carbonic acid, hexadecyl isobuty...	342 C21H42O3	1000314-61-3	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF022618\
 Data File : BF103232.D
 Acq On : 26 Feb 2018 21:57
 Operator : SJ/JU
 Sample : J1691-53 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B14

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF022218.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
unknown6.65	6.65	34.8	ng	1708490	1	6.87	981968	20.0
Hexadecane	10.33	2.3	ng	129467	3	9.92	1147740	20.0
Benzophenone	10.62	4.7	ng	270483	3	9.92	1147740	20.0
Heptacosane	11.25	2.3	ng	110537	4	11.40	976757	20.0
Anthracene, 1-met...	11.90	2.4	ng	118127	4	11.40	976757	20.0
1H-Cyclopropa[l]p...	11.93	2.5	ng	124012	4	11.40	976757	20.0
4H-Cyclopenta[def...	12.02	4.4	ng	213771	4	11.40	976757	20.0
unknown12.04	12.04	2.2	ng	106445	4	11.40	976757	20.0
18-Norabietane	12.26	2.2	ng	107816	4	11.40	976757	20.0
Fluoranthene, 2-m...	13.18	2.6	ng	146424	5	14.06	1106530	20.0
11H-Benzo[b]fluorene	13.24	2.7	ng	147005	5	14.06	1106530	20.0
1-Tricosene	13.87	2.8	ng	152684	5	14.06	1106530	20.0