

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF030618\
 Data File : BF103457.D
 Acq On : 6 Mar 2018 15:50
 Operator : SJ/JU
 Sample : J1724-19
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-12

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	5.104	510	513	524	rBV	57425	89485	2.56%	0.248%
2	5.498	576	580	607	rBV	2452107	3316276	94.69%	9.203%
3	6.516	748	753	771	rBV	2345545	3502213	100.00%	9.719%
4	6.645	771	775	797	rVB	2995222	3342718	95.45%	9.277%
5	6.869	810	813	821	rBV	783495	811948	23.18%	2.253%
6	7.028	836	840	856	rBV	2230582	2349995	67.10%	6.522%
7	7.439	906	910	924	rBV	1813059	2107484	60.18%	5.849%
8	8.157	1028	1032	1050	rVB	885032	1243832	35.52%	3.452%
9	8.728	1124	1129	1133	rBV	41265	40501	1.16%	0.112%
10	8.798	1138	1141	1150	rVB10	13147	35493	1.01%	0.099%
11	8.875	1150	1154	1160	rBV	54236	69390	1.98%	0.193%
12	9.228	1210	1214	1223	rBV	3127796	3010302	85.95%	8.354%
13	9.292	1223	1225	1229	rVV	110066	101930	2.91%	0.283%
14	9.333	1229	1232	1237	rVV	31654	37439	1.07%	0.104%
15	9.622	1275	1281	1287	rBV	187879	263187	7.51%	0.730%
16	9.775	1303	1307	1312	rBV	82436	101194	2.89%	0.281%
17	9.822	1312	1315	1319	rVV	179543	153546	4.38%	0.426%
18	9.910	1327	1330	1334	rBV	1039758	1036156	29.59%	2.876%
19	9.945	1334	1336	1339	rVB	66783	62286	1.78%	0.173%
20	10.057	1352	1355	1358	rBV3	29719	43721	1.25%	0.121%
21	10.122	1362	1366	1368	rVV2	61958	79989	2.28%	0.222%
22	10.145	1368	1370	1374	rVB3	48695	62926	1.80%	0.175%
23	10.227	1380	1384	1387	rBV4	28435	44130	1.26%	0.122%
24	10.322	1397	1400	1404	rVB	221001	186017	5.31%	0.516%
25	10.463	1421	1424	1431	rVB	71605	88926	2.54%	0.247%
26	10.545	1433	1438	1440	rBV2	96099	125954	3.60%	0.350%
27	10.622	1449	1451	1456	rVB4	36549	40753	1.16%	0.113%
28	10.710	1462	1466	1478	rBV	1478504	1680852	47.99%	4.665%
29	10.798	1478	1481	1491	rVB	250667	323322	9.23%	0.897%
30	10.951	1504	1507	1511	rVB	45873	44283	1.26%	0.123%
31	11.039	1520	1522	1527	rVB3	38378	40917	1.17%	0.114%
32	11.222	1550	1553	1555	rBV	38973	35777	1.02%	0.099%
33	11.245	1555	1557	1560	rVV	165705	159102	4.54%	0.442%
34	11.275	1560	1562	1565	rVV2	70050	71675	2.05%	0.199%

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35	11.304	1565	1567	1572	rVB	56611	51741	1.48%	0.144%
36	11.410	1581	1585	1587	rBV	922824	959380	27.39%	2.662%
37	11.433	1587	1589	1593	rVV	785614	666208	19.02%	1.849%
38	11.486	1595	1598	1602	rVV	183308	167513	4.78%	0.465%
39	11.645	1620	1625	1628	rBV2	76743	108403	3.10%	0.301%
40	11.674	1628	1630	1632	rVB	108643	75989	2.17%	0.211%
41	11.916	1667	1671	1673	rBV2	257886	289556	8.27%	0.804%
42	11.939	1673	1675	1681	rVB2	167254	163555	4.67%	0.454%
43	11.986	1681	1683	1686	rVB	56216	48149	1.37%	0.134%
44	12.022	1686	1689	1692	rBV	173785	207350	5.92%	0.575%
45	12.080	1697	1699	1703	rBV3	79103	78068	2.23%	0.217%
46	12.204	1717	1720	1724	rBV	92418	89843	2.57%	0.249%
47	12.245	1724	1727	1731	rVB2	68895	76229	2.18%	0.212%
48	12.463	1761	1764	1768	rBV2	88766	152625	4.36%	0.424%
49	12.557	1778	1780	1783	rBV	45250	43714	1.25%	0.121%
50	12.627	1788	1792	1796	rBV	1196735	1145432	32.71%	3.179%
51	12.863	1825	1832	1837	rBV	1049274	1290857	36.86%	3.582%
52	12.992	1850	1854	1858	rBV	2018242	1926875	55.02%	5.347%
53	13.074	1865	1868	1873	rBV3	59907	112614	3.22%	0.313%
54	13.186	1885	1887	1892	rVB2	162854	188049	5.37%	0.522%
55	13.257	1896	1899	1901	rBV	77517	74284	2.12%	0.206%
56	13.298	1903	1906	1909	rVB2	71321	75025	2.14%	0.208%
57	13.421	1925	1927	1932	rBV2	66944	76141	2.17%	0.211%
58	13.716	1974	1977	1984	rVB	93264	100064	2.86%	0.278%
59	13.886	2003	2006	2011	rBV4	365589	443383	12.66%	1.230%
60	14.086	2035	2040	2043	rBV2	732180	1153979	32.95%	3.203%
61	14.115	2043	2045	2048	rVB	436780	370583	10.58%	1.028%
62	15.192	2224	2228	2231	rBV	294691	449772	12.84%	1.248%
63	15.510	2279	2282	2286	rVB	159511	195965	5.60%	0.544%
64	15.574	2290	2293	2297	rVB	224593	253315	7.23%	0.703%
65	15.639	2301	2304	2307	rVB	259462	294913	8.42%	0.818%

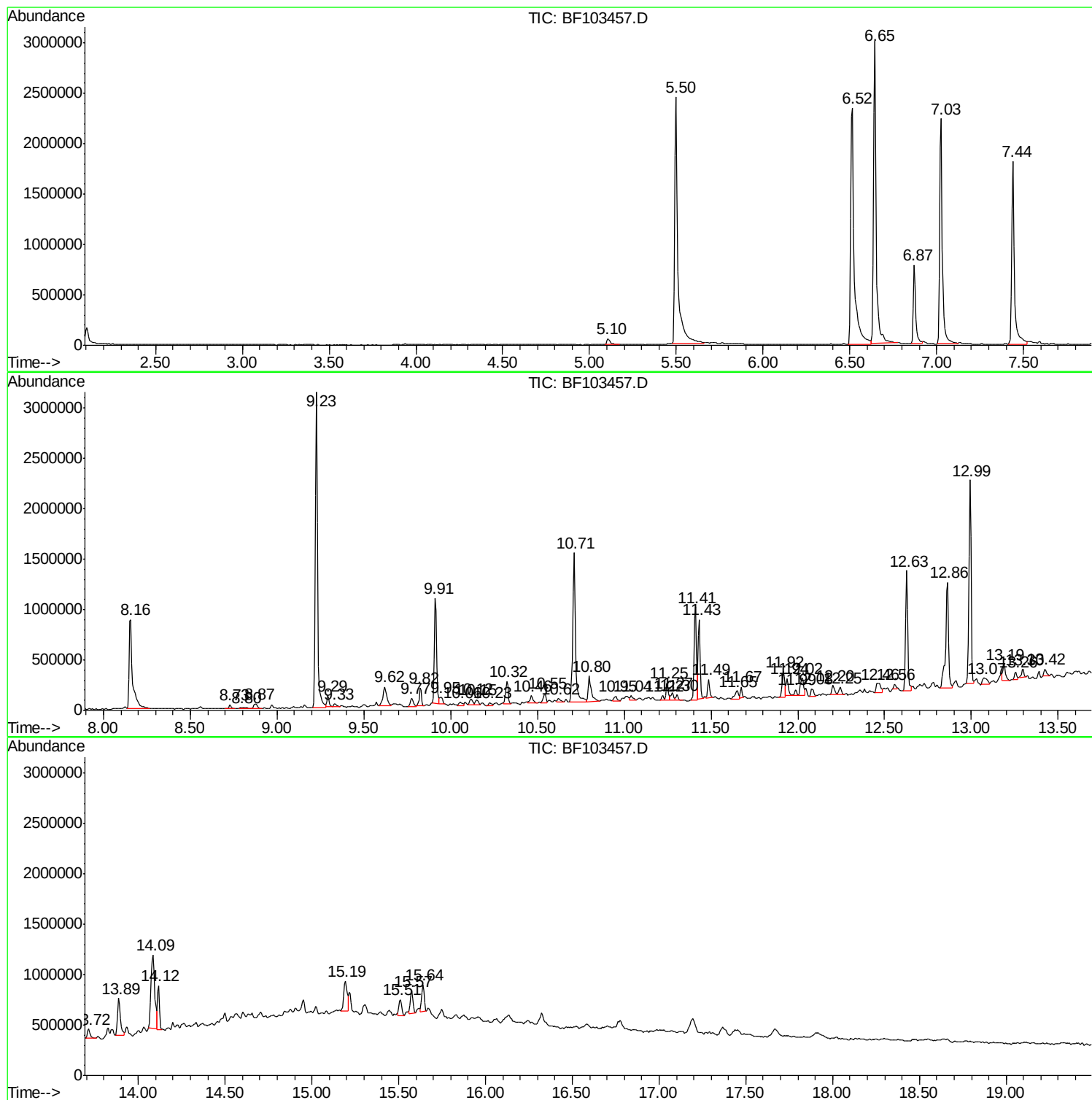
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
SP-12

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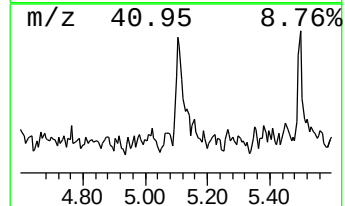
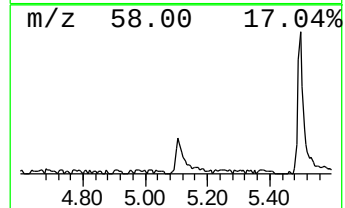
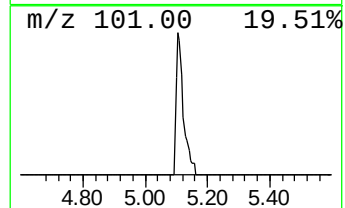
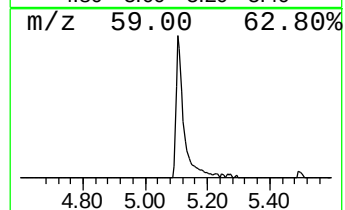
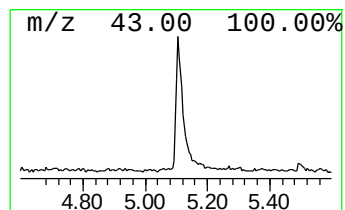
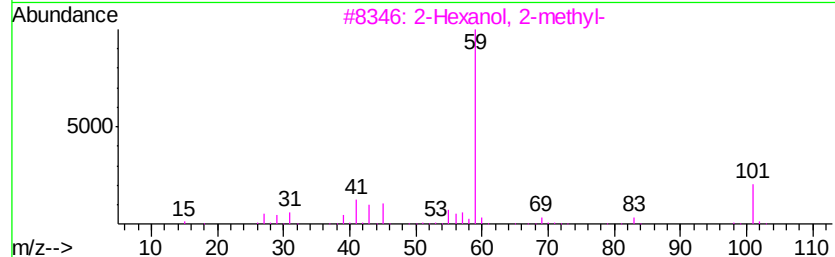
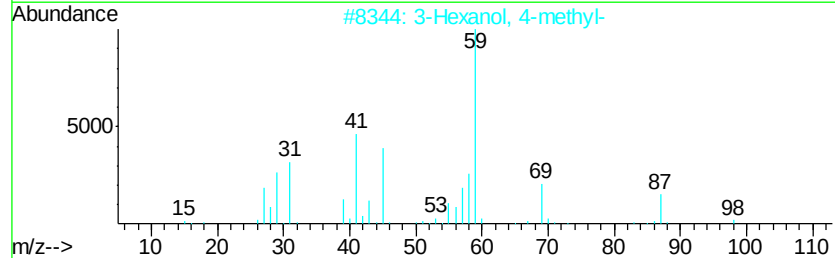
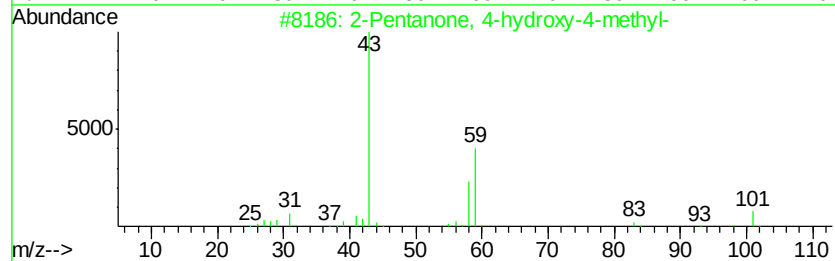
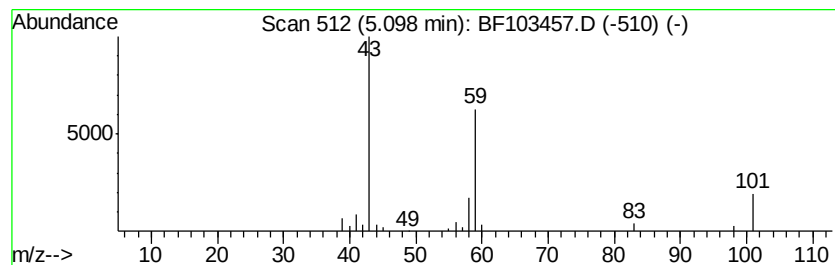
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	2.20 ng	89485	1,4-Dichlorobenzene-d4	6.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	38
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5			Guanidine	59	CH5N3	000113-00-8	9



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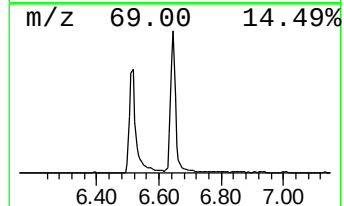
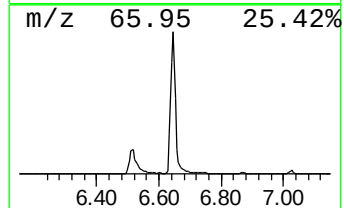
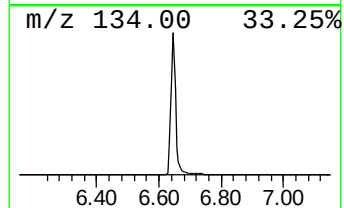
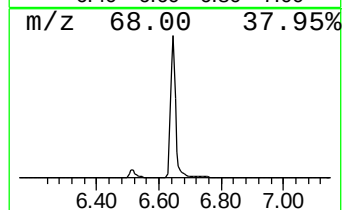
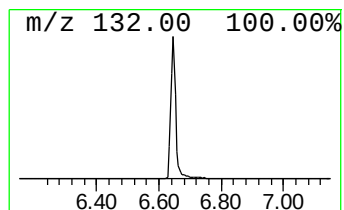
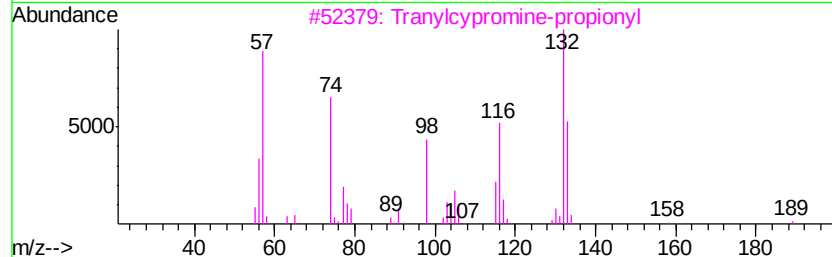
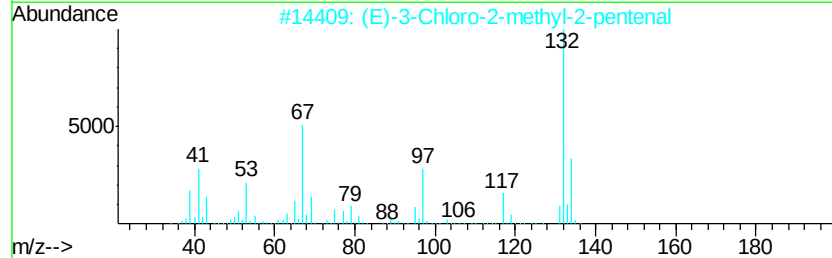
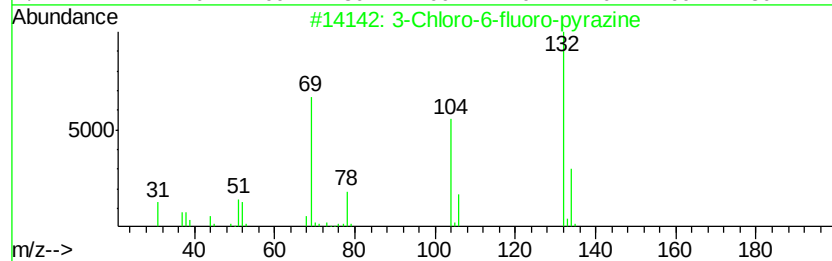
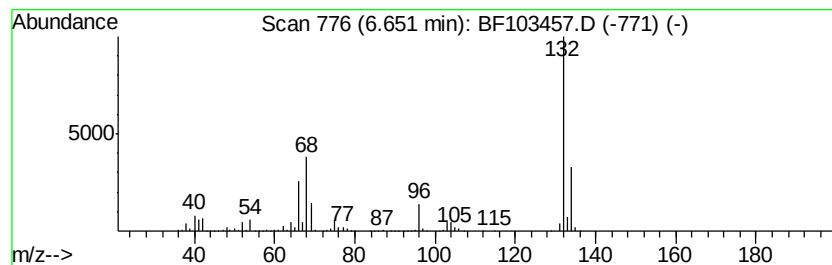
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	82.34 ng	3342720	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	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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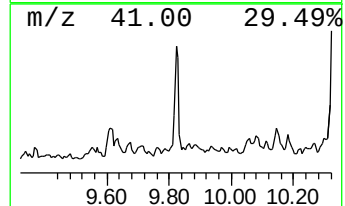
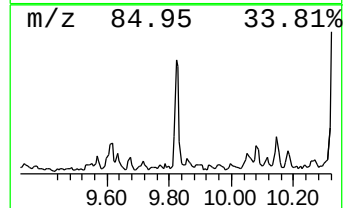
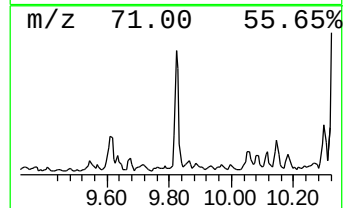
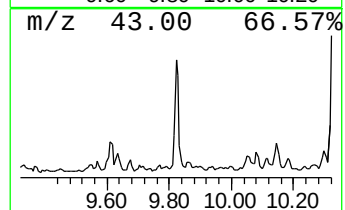
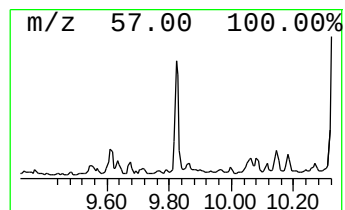
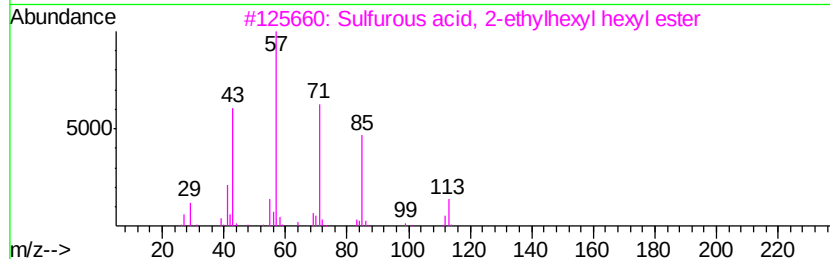
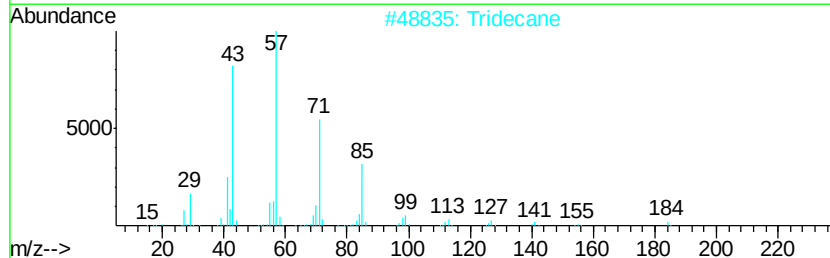
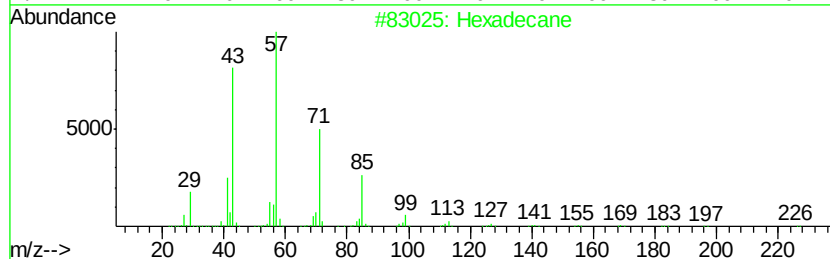
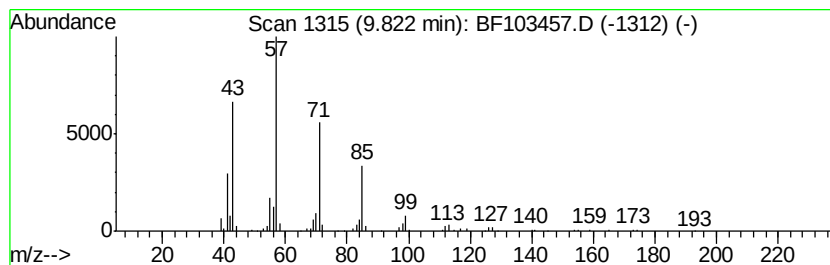
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	2.96 ng	153546	Acenaphthene-d10	9.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Tridecane	184	C13H28	000629-50-5	90
3		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	80
4		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	78
5		Undecane	156	C11H24	001120-21-4	64



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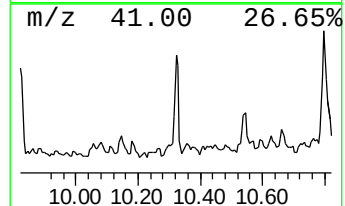
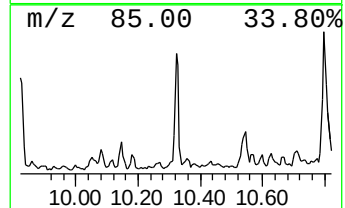
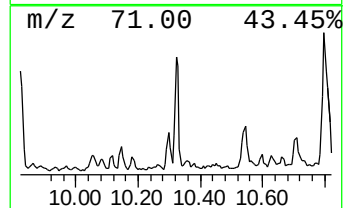
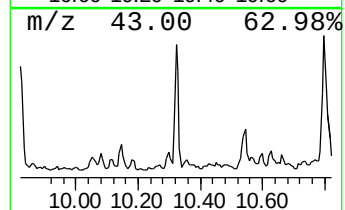
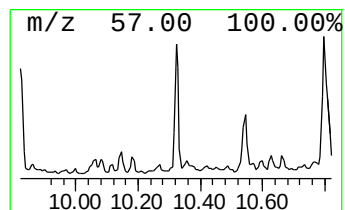
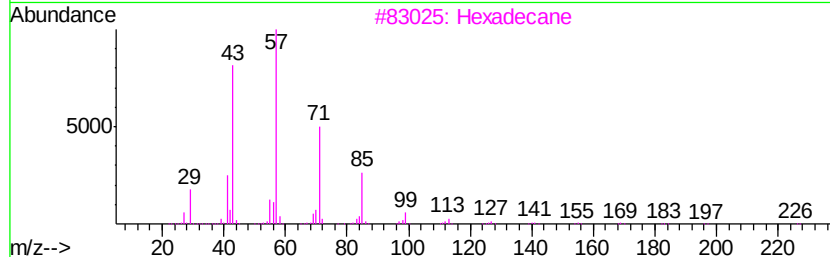
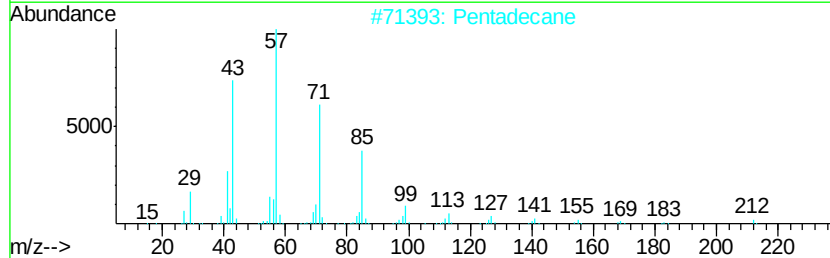
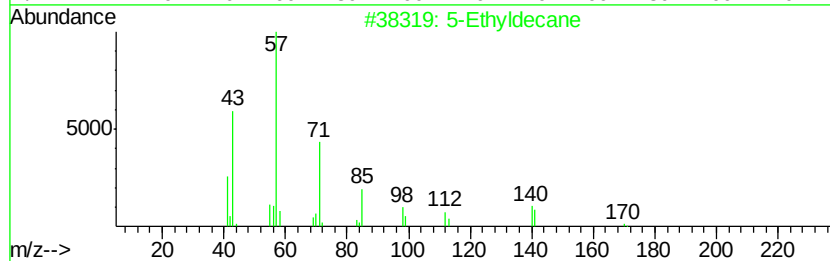
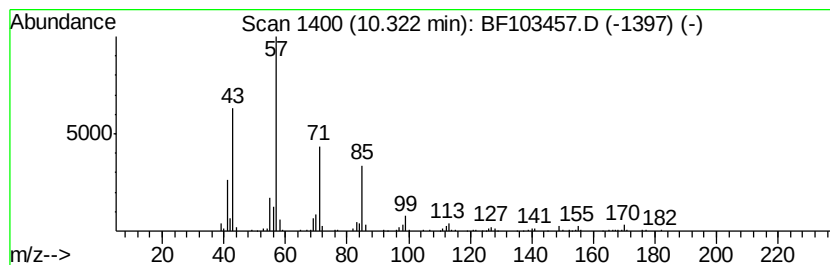
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 5-Ethyldecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	3.59 ng	186017	Acenaphthene-d10	9.91
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	5-Ethyldecane	170 C12H26	017302-36-2	91
2	Pentadecane	212 C15H32	000629-62-9	90
3	Hexadecane	226 C16H34	000544-76-3	86
4	Tetradecane	198 C14H30	000629-59-4	80
5	Dodecane, 2-methyl-6-propyl-	226 C16H34	055045-08-4	72



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ALS Vial : 13 Sample Multiplier: 1

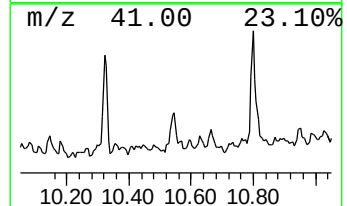
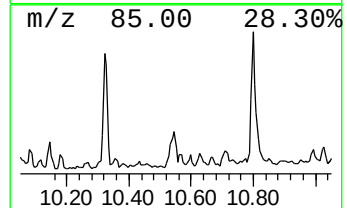
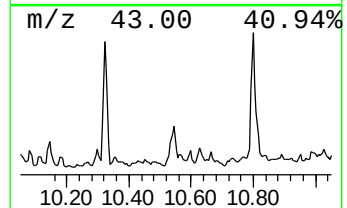
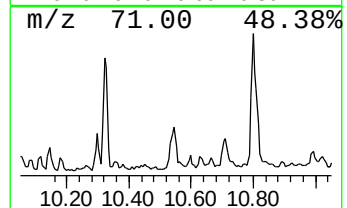
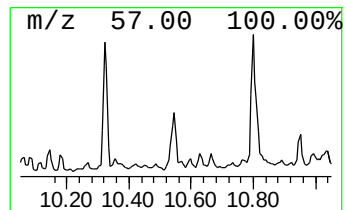
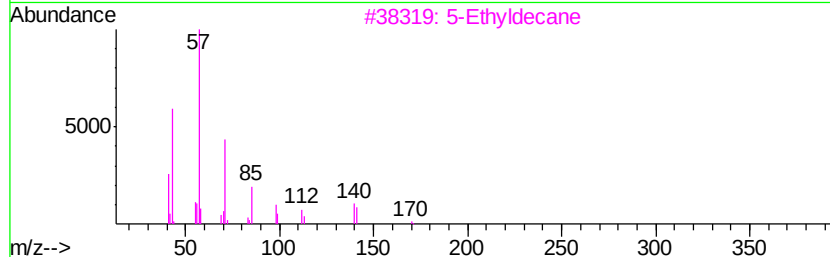
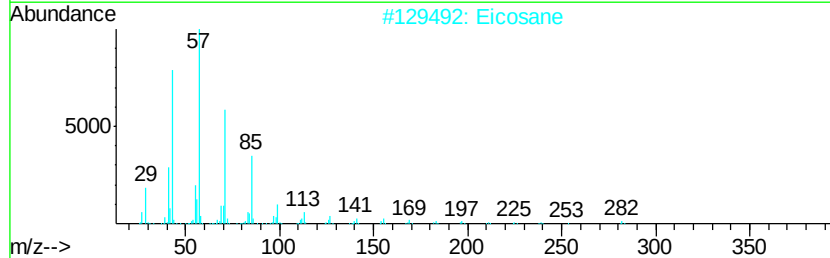
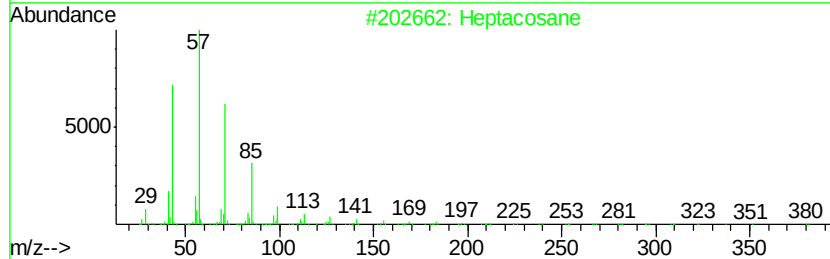
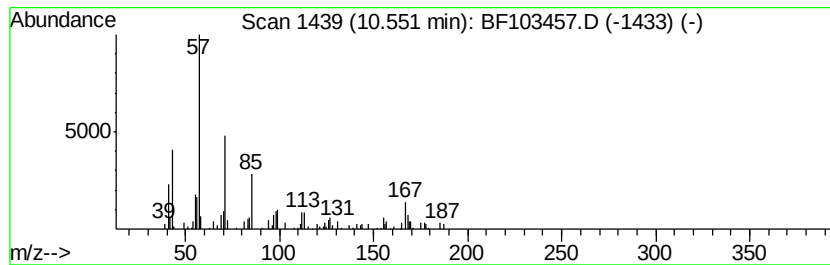
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heptacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.55	2.43 ng	125954	Acenaphthene-d10	9.91
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptacosane	380 C27H56	000593-49-7	80
2	Eicosane	282 C20H42	000112-95-8	80
3	5-Ethyldecane	170 C12H26	017302-36-2	74
4	Silane, trichlorooctadecyl-	386 C18H37Cl3Si	000112-04-9	72
5	Octacosane	394 C28H58	000630-02-4	64



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Operator : SJ/JU
Sample : J1724-19
Misc :
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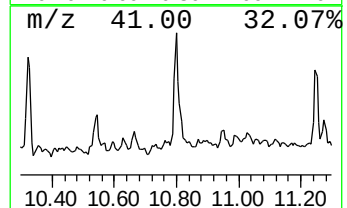
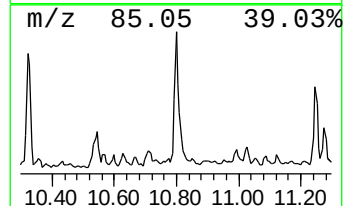
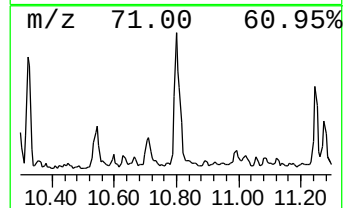
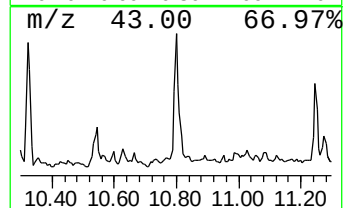
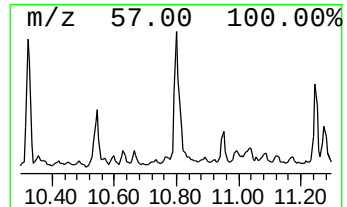
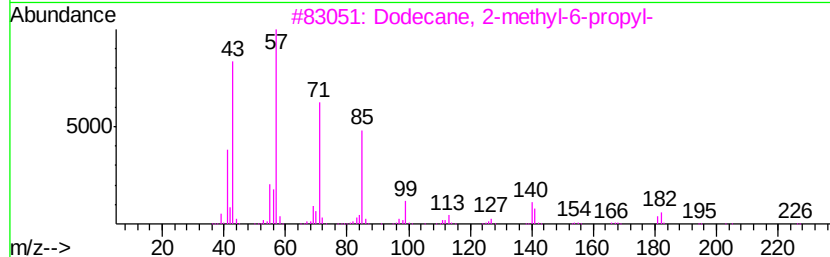
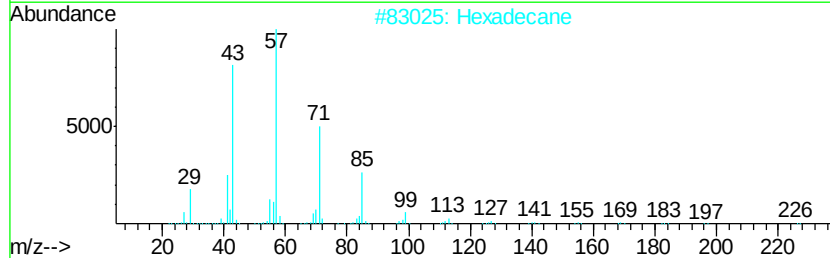
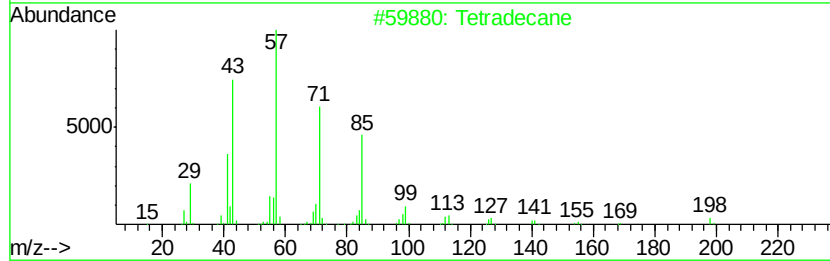
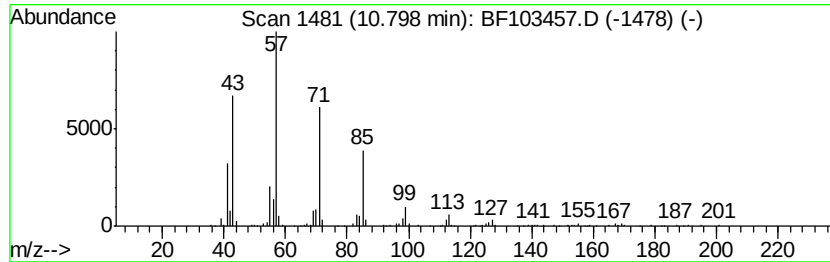
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetradecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.80	6.74 ng	323322	Phenanthrene-d10		11.41
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	90
2	Hexadecane	226	C16H34	000544-76-3	90
3	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
4	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	83
5	Undecane, 5-methyl-	170	C12H26	001632-70-8	72



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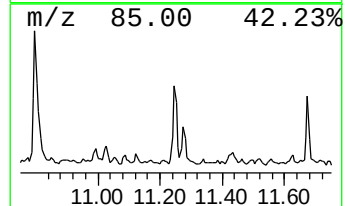
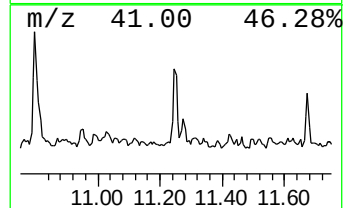
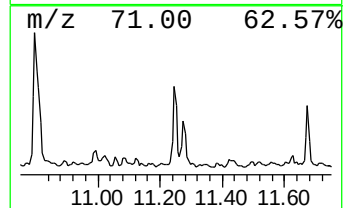
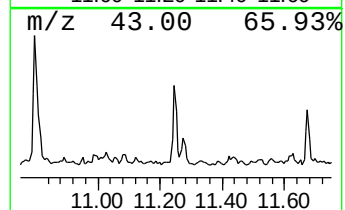
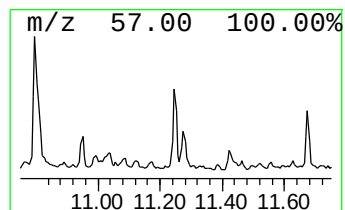
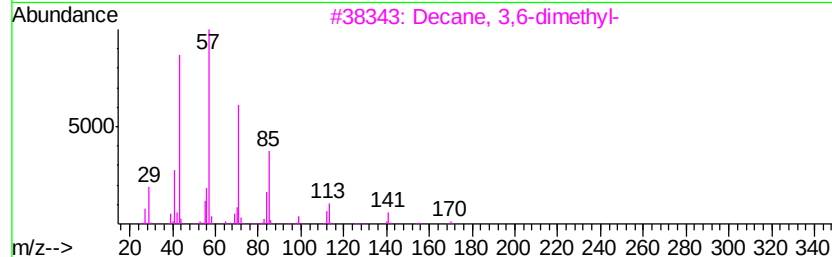
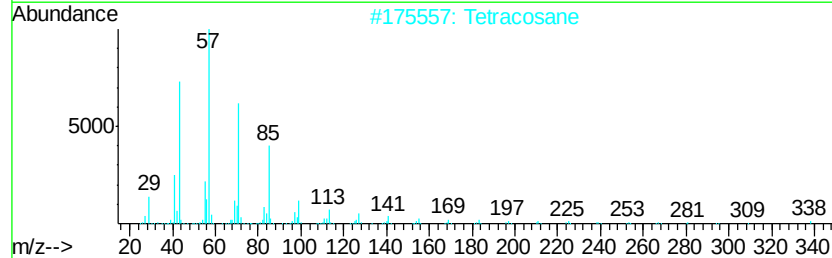
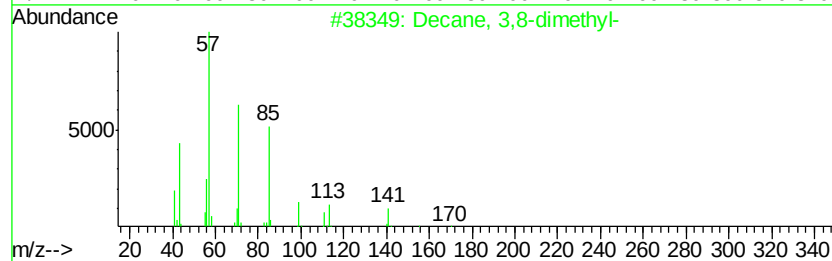
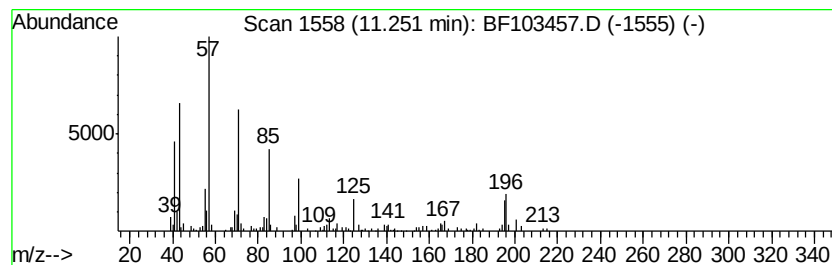
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Decane, 3,8-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.25	3.32 ng	159102	Phenanthrene-d10	11.41	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
2	Tetracosane	338	C24H50	000646-31-1	72
3	Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	68
4	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	64
5	Decane, 3-methyl-	156	C11H24	013151-34-3	64



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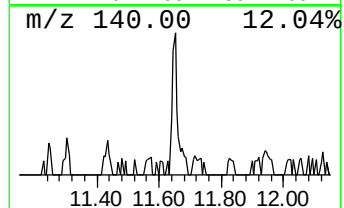
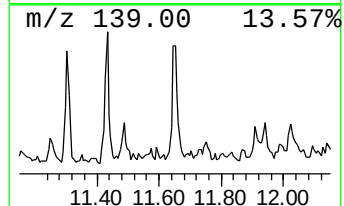
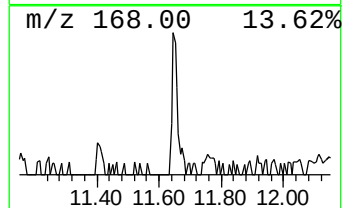
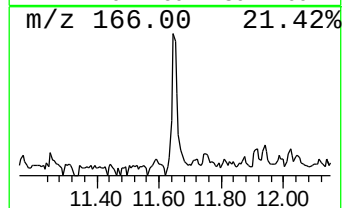
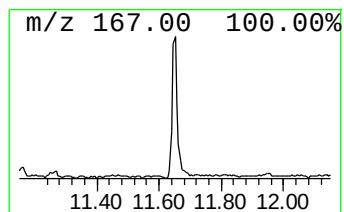
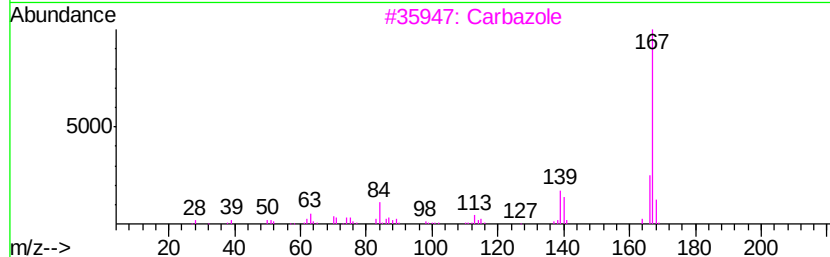
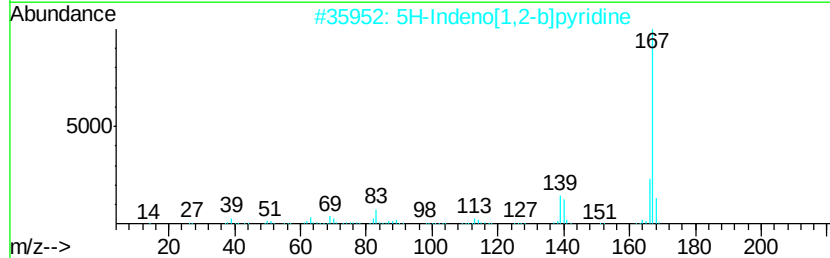
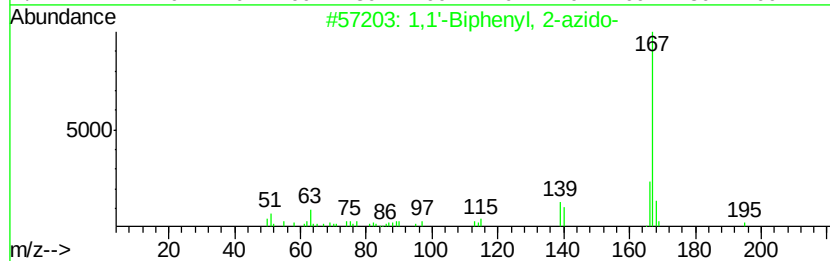
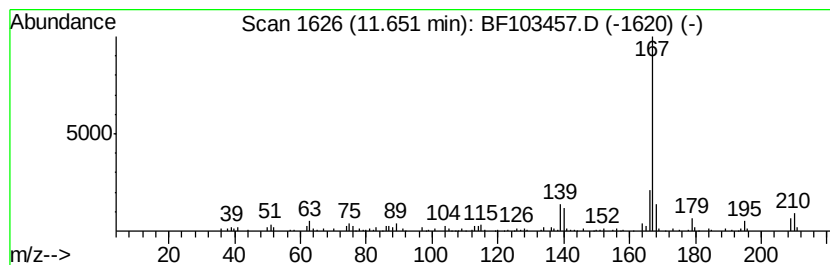
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1,1'-Biphenyl, 2-azido- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.65	2.26 ng	108403	Phenanthrene-d10	11.41	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-azido-	195	C12H9N3	007599-23-7	81
2	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	81
3	Carbazole	167	C12H9N	000086-74-8	81
4	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	72
5	9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	72



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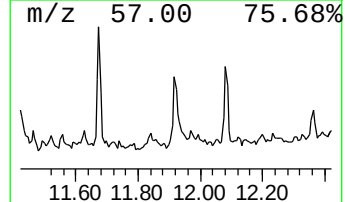
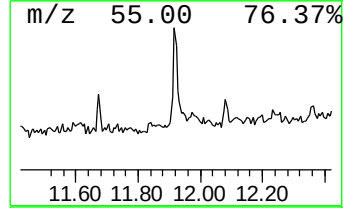
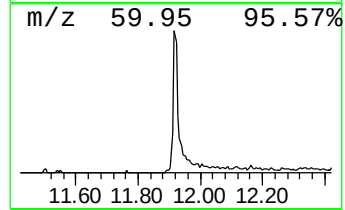
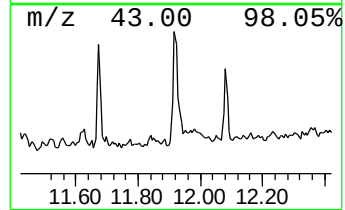
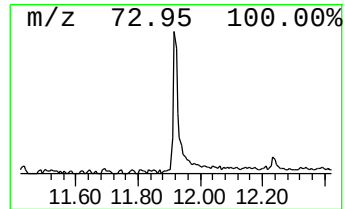
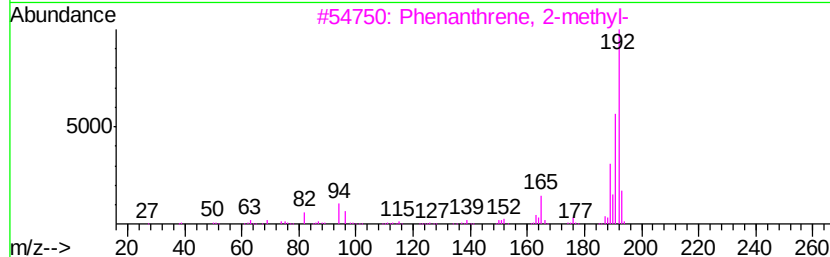
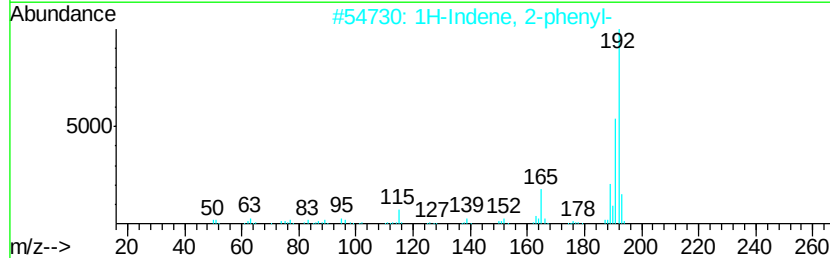
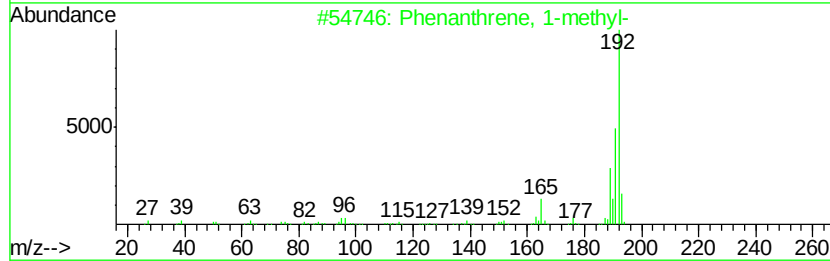
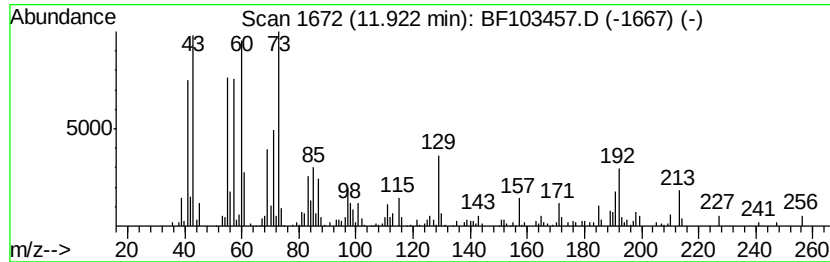
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.92	6.04 ng	289556	Phenanthrene-d10	11.41	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
2	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	91
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
5	Tridecanoic acid	214	C13H26O2	000638-53-9	87



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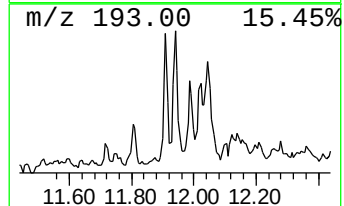
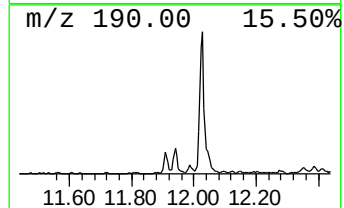
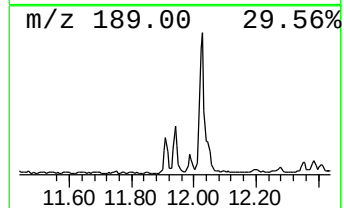
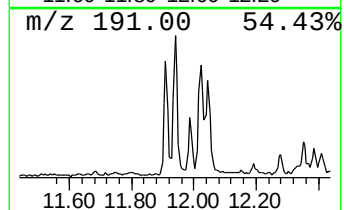
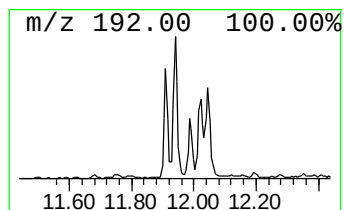
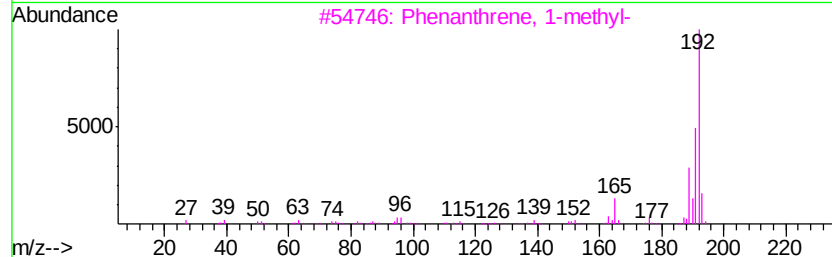
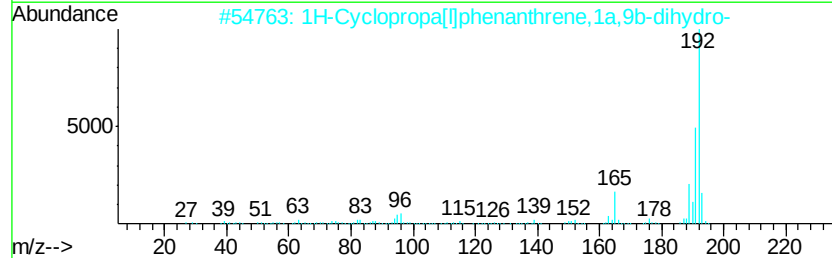
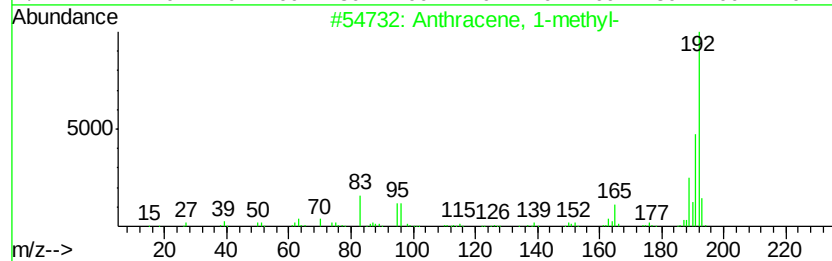
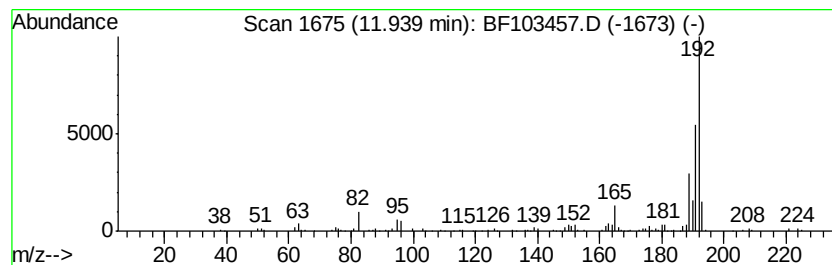
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Anthracene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	3.41 ng	163555	Phenanthrene-d10	11.41

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
2			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
5			1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	90



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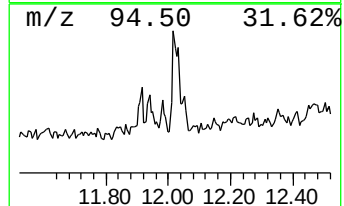
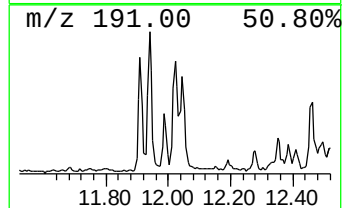
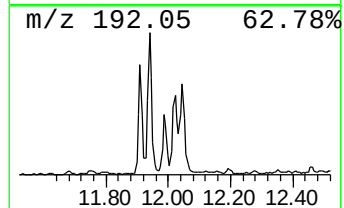
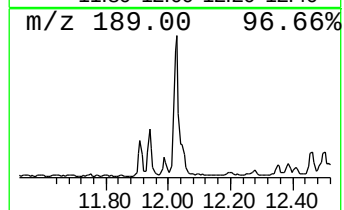
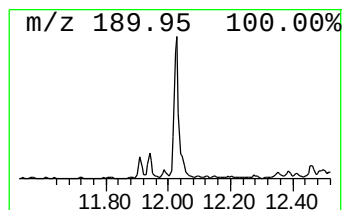
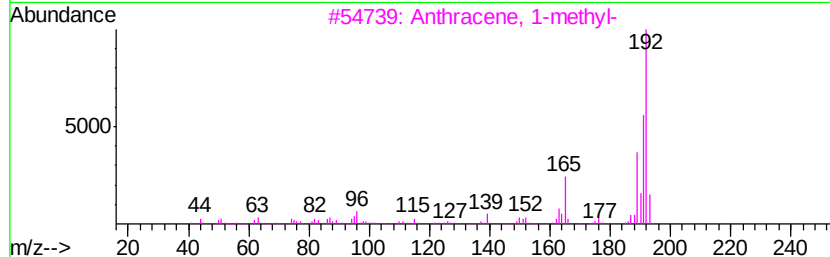
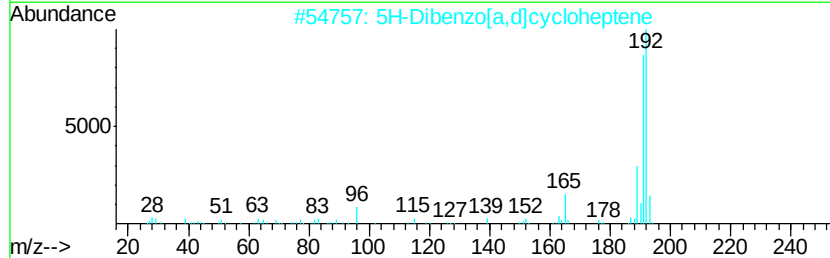
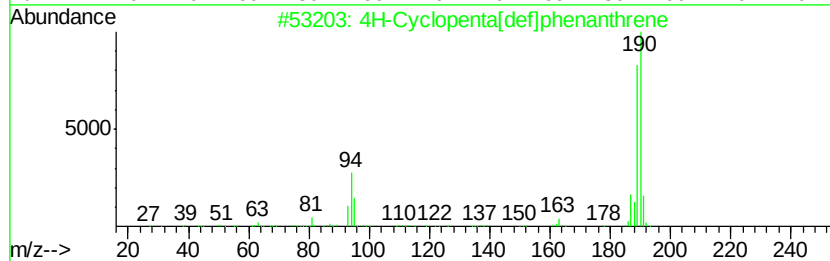
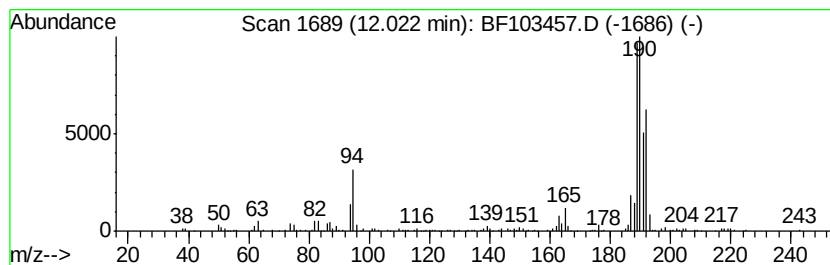
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	4.32 ng	207350	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	35
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	25
4		Benzene, 1,1'-(1,2-propadienyld...	192	C15H12	014251-57-1	18
5		Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	14



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Data File : BF103457.D
Acq On : 6 Mar 2018 15:50
Operator : SJ/JU
Sample : J1724-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

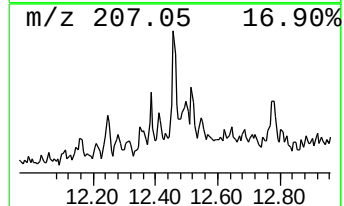
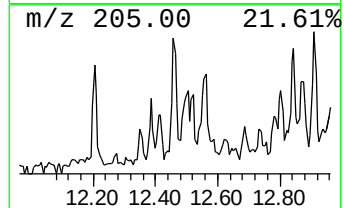
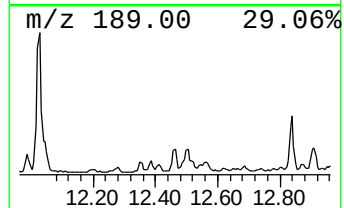
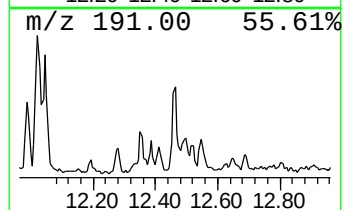
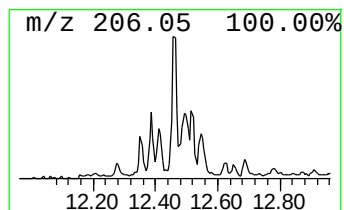
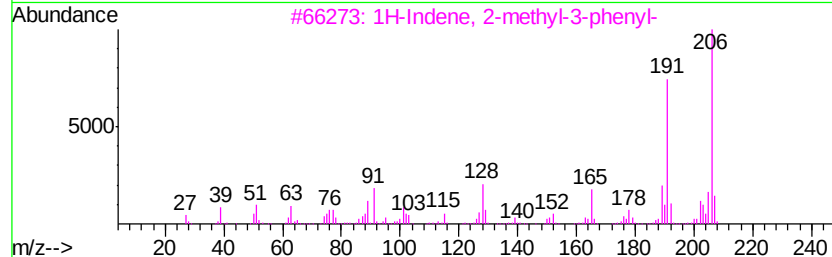
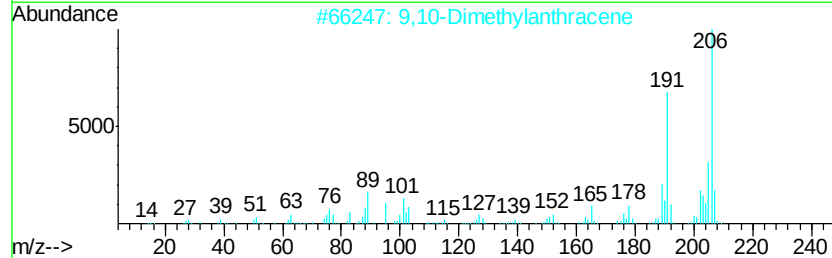
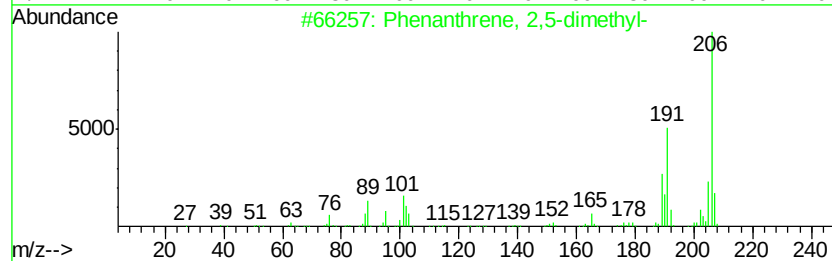
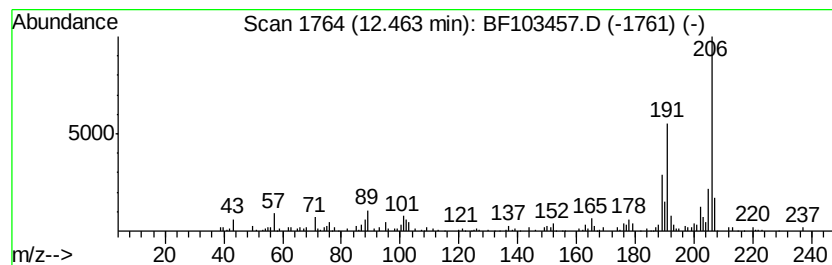
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phenanthrene, 2,5-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.46	3.18 ng	152625	Phenanthrene-d10		11.41
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2	9,10-Dimethylantracene	206	C16H14	000781-43-1	93
3	1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	91
4	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	89
5	Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	83



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF030618\
Data File : BF103457.D
Acq On : 6 Mar 2018 15:50
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Sample : J1724-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

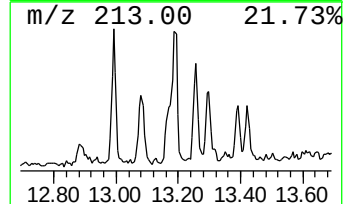
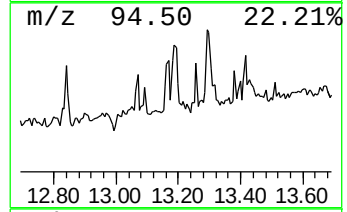
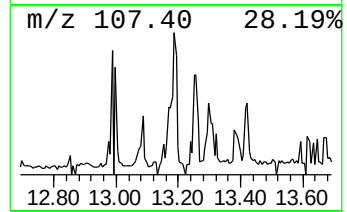
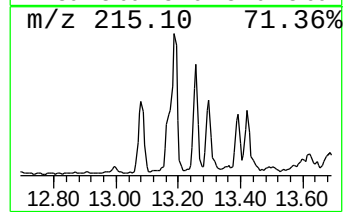
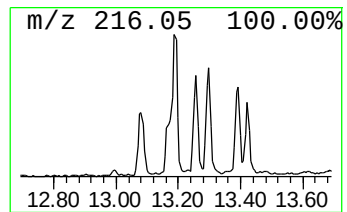
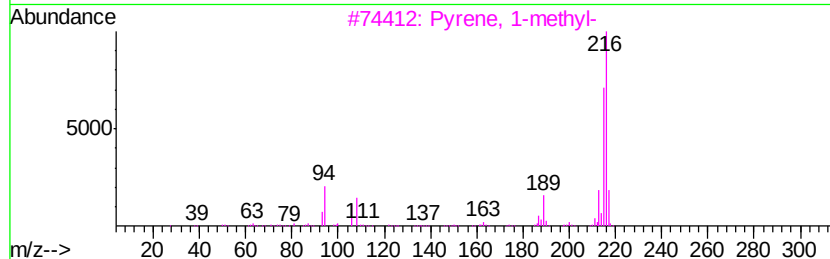
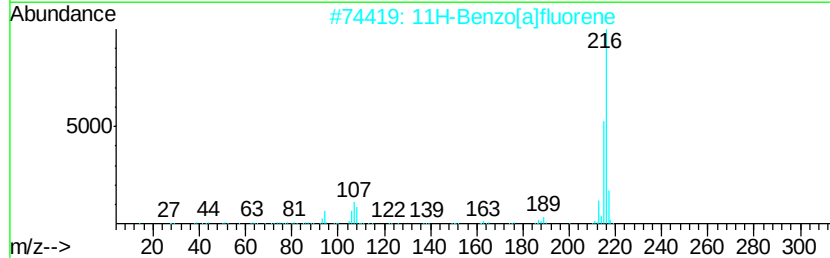
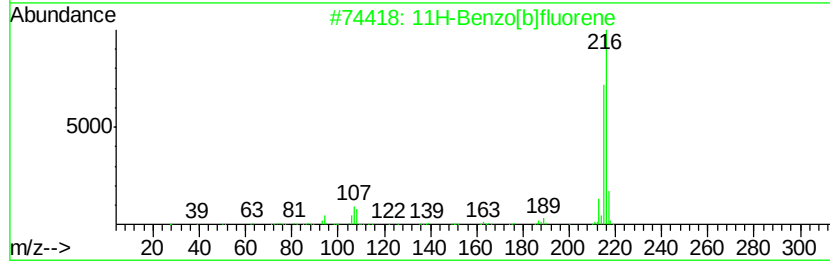
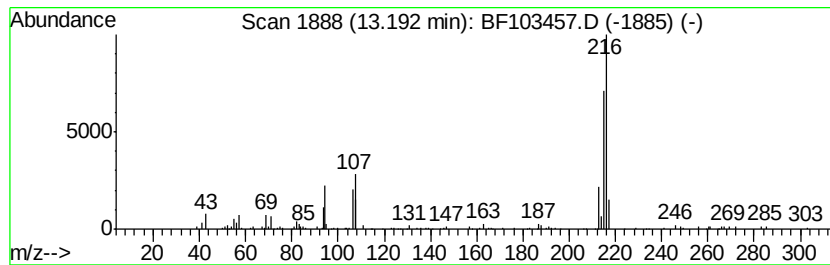
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 11H-Benzo[b]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.19	3.26 ng	188049	Chrysene-d12	14.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 76
2		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 68
3		Pvrene, 1-methyl-	216 C17H12	002381-21-7 58
4		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 50
5		Allopregnane-3.alpha.,20.alpha.-...	320 C21H36O2	000566-58-5 49



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030618\
Data File : BF103457.D
Acq On : 6 Mar 2018 15:50
Operator : SJ/JU
Sample : J1724-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP-12

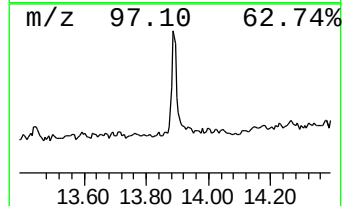
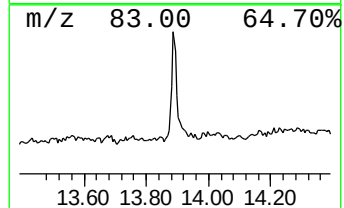
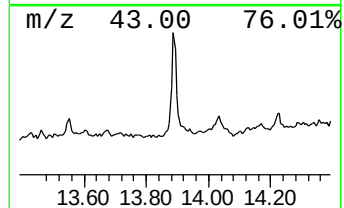
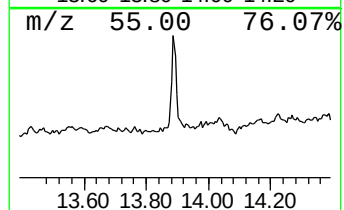
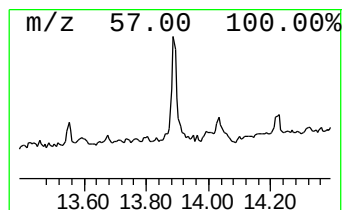
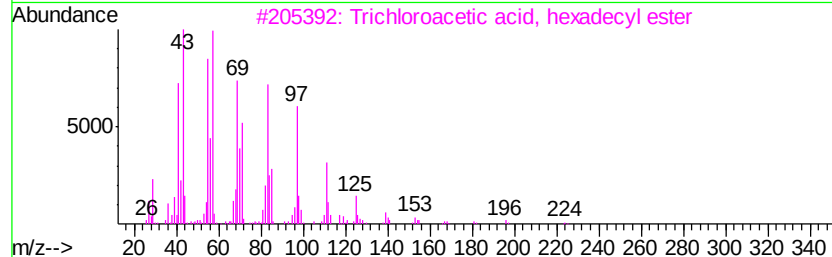
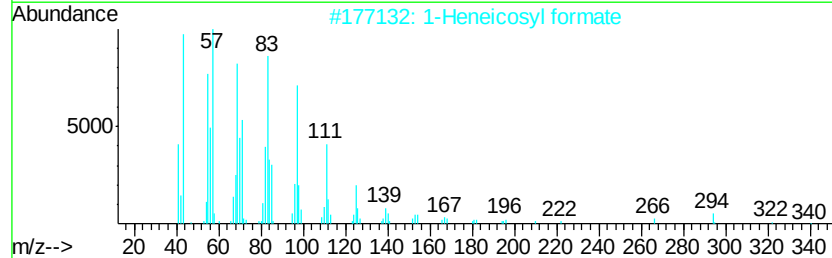
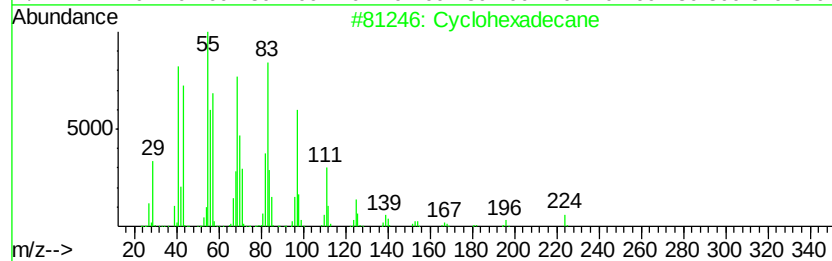
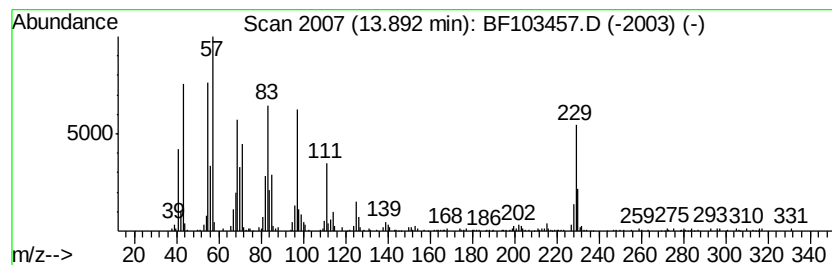
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Cyclohexadecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	7.68 ng	443383	Chrysene-d12	14.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	94
2		1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
4		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
5		Cyclotetracosane	336	C24H48	000297-03-0	91



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF030618\
Data File : BF103457.D
Acq On : 6 Mar 2018 15:50
Operator : SJ/JU
Sample : J1724-19
Misc :
ALS Vial : 13 Sample Multiplier: 1

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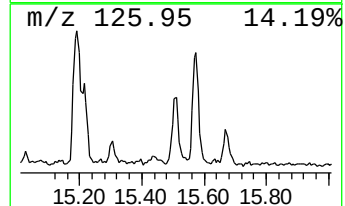
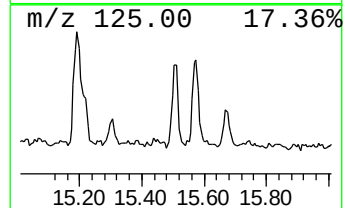
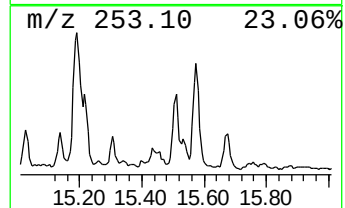
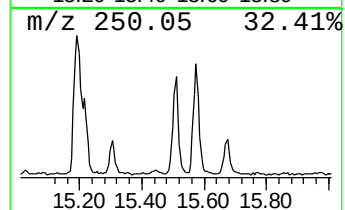
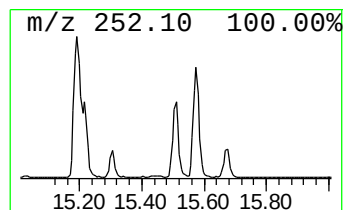
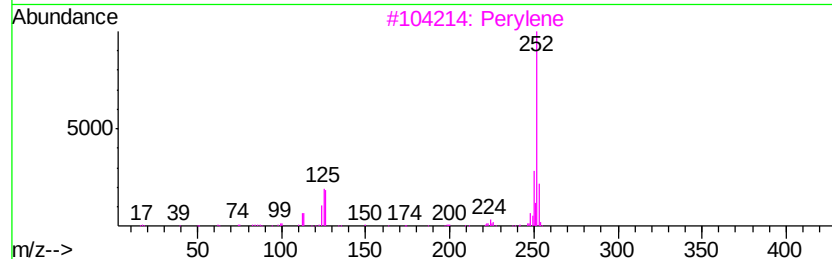
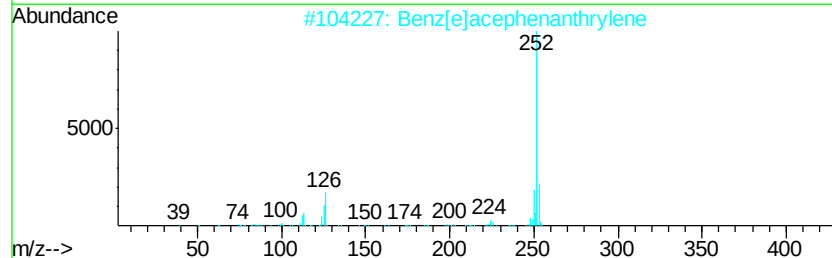
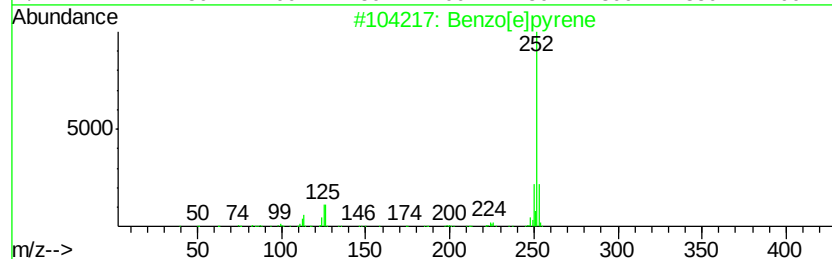
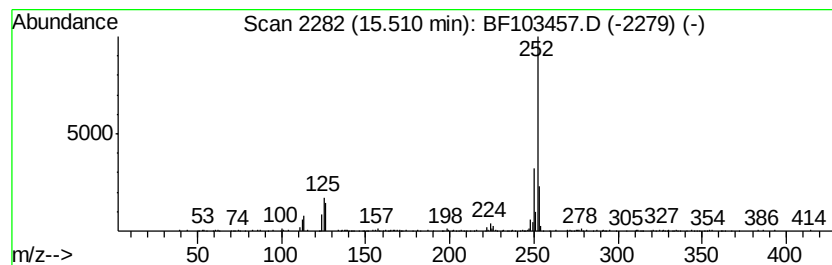
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.51	13.29 ng	195965	Perylene-d12	15.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	97
2		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	96
3		Perylene	252	C20H12	000198-55-0	94
4		Benzo[a]pyrene	252	C20H12	000050-32-8	90
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF030618\
 Data File : BF103457.D
 Acq On : 6 Mar 2018 15:50
 Operator : SJ/JU
 Sample : J1724-19
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	2.2	ng	89485	1	6.87	811948	20.0
unknown6.65	6.65	82.3	ng	3342720	1	6.87	811948	20.0
Hexadecane	9.82	3.0	ng	153546	3	9.91	1036160	20.0
5-Ethyldecane	10.32	3.6	ng	186017	3	9.91	1036160	20.0
Heptacosane	10.55	2.4	ng	125954	3	9.91	1036160	20.0
Tetradecane	10.80	6.7	ng	323322	4	11.41	959380	20.0
Decane, 3,8-dimet...	11.25	3.3	ng	159102	4	11.41	959380	20.0
1,1'-Biphenyl, 2-...	11.65	2.3	ng	108403	4	11.41	959380	20.0
Phenanthrene, 1-m...	11.92	6.0	ng	289556	4	11.41	959380	20.0
Anthracene, 1-met...	11.94	3.4	ng	163555	4	11.41	959380	20.0
4H-Cyclopenta[def...	12.02	4.3	ng	207350	4	11.41	959380	20.0
Phenanthrene, 2,5...	12.46	3.2	ng	152625	4	11.41	959380	20.0
11H-Benzo[b]fluorene	13.19	3.3	ng	188049	5	14.09	1153980	20.0
Cyclohexadecane	13.89	7.7	ng	443383	5	14.09	1153980	20.0
Benzo[e]pyrene	15.51	13.3	ng	195965	6	15.64	294913	20.0