

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
 Data File : BF097443.D  
 Acq On : 7 Aug 2017 21:48  
 Operator : SJ/JU  
 Sample : I4623-06  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TEC-WN

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-BF072517.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.534       | 464           | 467         | 478          | rBV      | 129207         | 211365        | 9.80%           | 0.679%        |
| 2         | 5.004       | 543           | 547         | 568          | rBV      | 1537230        | 1967013       | 91.22%          | 6.323%        |
| 3         | 6.081       | 727           | 730         | 744          | rBV      | 1868944        | 2156308       | 100.00%         | 6.932%        |
| 4         | 6.187       | 744           | 748         | 757          | rVB      | 1867660        | 1942854       | 90.10%          | 6.246%        |
| 5         | 6.410       | 783           | 786         | 797          | rBV      | 623520         | 570308        | 26.45%          | 1.833%        |
| 6         | 6.563       | 809           | 812         | 820          | rBV      | 1351810        | 1336807       | 62.00%          | 4.297%        |
| 7         | 6.981       | 880           | 883         | 897          | rBV      | 1319646        | 1323799       | 61.39%          | 4.256%        |
| 8         | 7.145       | 908           | 911         | 920          | rVB      | 31206          | 43123         | 2.00%           | 0.139%        |
| 9         | 7.692       | 1001          | 1004        | 1016         | rVB      | 856280         | 791306        | 36.70%          | 2.544%        |
| 10        | 8.410       | 1123          | 1126        | 1134         | rVB      | 31662          | 29753         | 1.38%           | 0.096%        |
| 11        | 8.769       | 1183          | 1187        | 1191         | rBV      | 2042709        | 1940775       | 90.00%          | 6.239%        |
| 12        | 9.175       | 1252          | 1256        | 1262         | rBV      | 559986         | 464008        | 21.52%          | 1.492%        |
| 13        | 9.433       | 1297          | 1300        | 1303         | rBV      | 831989         | 751220        | 34.84%          | 2.415%        |
| 14        | 9.469       | 1303          | 1306        | 1314         | rVB      | 349377         | 301763        | 13.99%          | 0.970%        |
| 15        | 9.645       | 1333          | 1336        | 1341         | rVB      | 86333          | 79784         | 3.70%           | 0.256%        |
| 16        | 9.892       | 1375          | 1378        | 1381         | rVB      | 52496          | 42703         | 1.98%           | 0.137%        |
| 17        | 9.980       | 1390          | 1393        | 1400         | rBV      | 173832         | 145531        | 6.75%           | 0.468%        |
| 18        | 10.045      | 1400          | 1404        | 1406         | rBV      | 22978          | 30072         | 1.39%           | 0.097%        |
| 19        | 10.110      | 1410          | 1415        | 1418         | rBV4     | 33332          | 46292         | 2.15%           | 0.149%        |
| 20        | 10.222      | 1431          | 1434        | 1440         | rBV      | 902088         | 862502        | 40.00%          | 2.773%        |
| 21        | 10.363      | 1455          | 1458        | 1466         | rBV      | 98266          | 120363        | 5.58%           | 0.387%        |
| 22        | 10.727      | 1517          | 1520        | 1524         | rBV      | 42420          | 42081         | 1.95%           | 0.135%        |
| 23        | 10.804      | 1530          | 1533        | 1536         | rBV      | 111753         | 100335        | 4.65%           | 0.323%        |
| 24        | 10.910      | 1547          | 1551        | 1553         | rBV      | 630190         | 598652        | 27.76%          | 1.924%        |
| 25        | 10.933      | 1553          | 1555        | 1559         | rVV      | 1291678        | 1030472       | 47.79%          | 3.313%        |
| 26        | 10.986      | 1559          | 1564        | 1569         | rVV      | 334568         | 314543        | 14.59%          | 1.011%        |
| 27        | 11.151      | 1586          | 1592        | 1595         | rBV2     | 145165         | 164196        | 7.61%           | 0.528%        |
| 28        | 11.233      | 1603          | 1606        | 1609         | rVB2     | 32272          | 35987         | 1.67%           | 0.116%        |
| 29        | 11.410      | 1632          | 1636        | 1638         | rBV      | 111237         | 89744         | 4.16%           | 0.288%        |
| 30        | 11.439      | 1638          | 1641        | 1644         | rVV      | 153670         | 142425        | 6.61%           | 0.458%        |
| 31        | 11.469      | 1644          | 1646        | 1647         | rVV      | 75173          | 57193         | 2.65%           | 0.184%        |
| 32        | 11.486      | 1647          | 1649        | 1652         | rVV      | 77437          | 71863         | 3.33%           | 0.231%        |
| 33        | 11.516      | 1652          | 1654        | 1657         | rVV      | 227628         | 218762        | 10.15%          | 0.703%        |
| 34        | 11.539      | 1657          | 1658        | 1665         | rVB      | 64896          | 61963         | 2.87%           | 0.199%        |

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Max Peaks: 100

Peak Location: TOP

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Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 11.633 | 1671 | 1674 | 1680 | rBV6 | 24987   | 39400   | 1.83%  | 0.127% |
| 36 | 11.704 | 1683 | 1686 | 1690 | rVV  | 86590   | 71201   | 3.30%  | 0.229% |
| 37 | 11.745 | 1690 | 1693 | 1696 | rVV  | 83509   | 80749   | 3.74%  | 0.260% |
| 38 | 11.957 | 1726 | 1729 | 1732 | rBV  | 41897   | 45551   | 2.11%  | 0.146% |
| 39 | 12.051 | 1742 | 1745 | 1748 | rVV2 | 55596   | 50366   | 2.34%  | 0.162% |
| 40 | 12.116 | 1752 | 1756 | 1760 | rBV  | 1517713 | 1353938 | 62.79% | 4.352% |
| 41 | 12.263 | 1776 | 1781 | 1784 | rBV3 | 65682   | 85194   | 3.95%  | 0.274% |
| 42 | 12.339 | 1789 | 1794 | 1797 | rBV  | 1225070 | 1289268 | 59.79% | 4.145% |
| 43 | 12.392 | 1800 | 1803 | 1811 | rVB2 | 54204   | 74254   | 3.44%  | 0.239% |
| 44 | 12.463 | 1811 | 1815 | 1816 | rBV  | 77062   | 61204   | 2.84%  | 0.197% |
| 45 | 12.492 | 1816 | 1820 | 1827 | rVB  | 1349067 | 1363814 | 63.25% | 4.384% |
| 46 | 12.568 | 1827 | 1833 | 1835 | rBV2 | 62420   | 103666  | 4.81%  | 0.333% |
| 47 | 12.598 | 1835 | 1838 | 1843 | rVB  | 42044   | 39624   | 1.84%  | 0.127% |
| 48 | 12.645 | 1843 | 1846 | 1848 | rBV3 | 54375   | 71865   | 3.33%  | 0.231% |
| 49 | 12.668 | 1848 | 1850 | 1854 | rVB  | 175307  | 167749  | 7.78%  | 0.539% |
| 50 | 12.733 | 1858 | 1861 | 1864 | rVB  | 122656  | 113123  | 5.25%  | 0.364% |
| 51 | 12.768 | 1864 | 1867 | 1869 | rBV  | 108925  | 115399  | 5.35%  | 0.371% |
| 52 | 12.863 | 1880 | 1883 | 1886 | rBV  | 60368   | 44671   | 2.07%  | 0.144% |
| 53 | 12.892 | 1886 | 1888 | 1892 | rBV  | 49372   | 43242   | 2.01%  | 0.139% |
| 54 | 13.074 | 1910 | 1919 | 1920 | rBV6 | 25402   | 53198   | 2.47%  | 0.171% |
| 55 | 13.186 | 1935 | 1938 | 1940 | rVV  | 61816   | 54354   | 2.52%  | 0.175% |
| 56 | 13.286 | 1951 | 1955 | 1956 | rBV  | 92783   | 95146   | 4.41%  | 0.306% |
| 57 | 13.304 | 1956 | 1958 | 1961 | rVV  | 110355  | 108225  | 5.02%  | 0.348% |
| 58 | 13.333 | 1961 | 1963 | 1965 | rVV  | 98174   | 92353   | 4.28%  | 0.297% |
| 59 | 13.351 | 1965 | 1966 | 1970 | rVB2 | 55661   | 50932   | 2.36%  | 0.164% |
| 60 | 13.392 | 1970 | 1973 | 1976 | rBV  | 88322   | 84646   | 3.93%  | 0.272% |
| 61 | 13.427 | 1976 | 1979 | 1989 | rVB2 | 54597   | 153263  | 7.11%  | 0.493% |
| 62 | 13.527 | 1989 | 1996 | 1999 | rBV2 | 901551  | 1407627 | 65.28% | 4.525% |
| 63 | 13.557 | 1999 | 2001 | 2005 | rVB  | 630954  | 593252  | 27.51% | 1.907% |
| 64 | 13.633 | 2010 | 2014 | 2019 | rVB  | 148095  | 145444  | 6.75%  | 0.468% |
| 65 | 13.692 | 2019 | 2024 | 2027 | rBV2 | 55818   | 76711   | 3.56%  | 0.247% |
| 66 | 13.733 | 2028 | 2031 | 2035 | rVV2 | 68124   | 96928   | 4.50%  | 0.312% |
| 67 | 13.768 | 2035 | 2037 | 2040 | rVB  | 55855   | 44466   | 2.06%  | 0.143% |
| 68 | 13.886 | 2054 | 2057 | 2059 | rVV2 | 46504   | 52768   | 2.45%  | 0.170% |
| 69 | 13.921 | 2059 | 2063 | 2066 | rVV  | 130945  | 156538  | 7.26%  | 0.503% |
| 70 | 13.951 | 2066 | 2068 | 2071 | rVV2 | 60182   | 61147   | 2.84%  | 0.197% |
| 71 | 14.015 | 2071 | 2079 | 2081 | rVV3 | 64378   | 125810  | 5.83%  | 0.404% |

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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

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Peak Location: TOP

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

Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |        |        |        |        |
|-----|--------|------|------|------|------|--------|--------|--------|--------|
| 72  | 14.045 | 2081 | 2084 | 2085 | rVV3 | 71385  | 81347  | 3.77%  | 0.262% |
| 73  | 14.063 | 2085 | 2087 | 2090 | rVV2 | 102385 | 90936  | 4.22%  | 0.292% |
| 74  | 14.245 | 2113 | 2118 | 2125 | rBV4 | 63079  | 90246  | 4.19%  | 0.290% |
| 75  | 14.321 | 2129 | 2131 | 2137 | rVB4 | 25895  | 30836  | 1.43%  | 0.099% |
| 76  | 14.398 | 2139 | 2144 | 2147 | rVB2 | 66161  | 72543  | 3.36%  | 0.233% |
| 77  | 14.486 | 2155 | 2159 | 2161 | rBV2 | 32782  | 37312  | 1.73%  | 0.120% |
| 78  | 14.527 | 2162 | 2166 | 2174 | rVV  | 592172 | 952513 | 44.17% | 3.062% |
| 79  | 14.592 | 2174 | 2177 | 2178 | rVV  | 30810  | 31347  | 1.45%  | 0.101% |
| 80  | 14.615 | 2178 | 2181 | 2185 | rVB  | 100128 | 101046 | 4.69%  | 0.325% |
| 81  | 14.698 | 2189 | 2195 | 2197 | rBV5 | 41381  | 80672  | 3.74%  | 0.259% |
| 82  | 14.727 | 2197 | 2200 | 2203 | rVV2 | 47382  | 75492  | 3.50%  | 0.243% |
| 83  | 14.768 | 2203 | 2207 | 2213 | rVV  | 341364 | 397740 | 18.45% | 1.279% |
| 84  | 14.821 | 2213 | 2216 | 2221 | rVV  | 486076 | 537575 | 24.93% | 1.728% |
| 85  | 14.874 | 2221 | 2225 | 2227 | rVV  | 449940 | 490858 | 22.76% | 1.578% |
| 86  | 14.898 | 2227 | 2229 | 2236 | rVB2 | 144845 | 190796 | 8.85%  | 0.613% |
| 87  | 15.062 | 2253 | 2257 | 2261 | rVB3 | 32528  | 47872  | 2.22%  | 0.154% |
| 88  | 15.251 | 2285 | 2289 | 2294 | rBV4 | 26324  | 41287  | 1.91%  | 0.133% |
| 89  | 15.333 | 2299 | 2303 | 2311 | rVB4 | 43600  | 75043  | 3.48%  | 0.241% |
| 90  | 15.651 | 2353 | 2357 | 2361 | rBV  | 24714  | 35551  | 1.65%  | 0.114% |
| 91  | 15.904 | 2394 | 2400 | 2402 | rVV2 | 24303  | 45195  | 2.10%  | 0.145% |
| 92  | 15.939 | 2403 | 2406 | 2409 | rVV  | 27194  | 34761  | 1.61%  | 0.112% |
| 93  | 15.980 | 2411 | 2413 | 2421 | rVB8 | 20023  | 41150  | 1.91%  | 0.132% |
| 94  | 16.062 | 2421 | 2427 | 2433 | rBV  | 139061 | 266243 | 12.35% | 0.856% |
| 95  | 16.115 | 2433 | 2436 | 2441 | rVB6 | 31549  | 47836  | 2.22%  | 0.154% |
| 96  | 16.198 | 2446 | 2450 | 2454 | rBV2 | 34071  | 60808  | 2.82%  | 0.195% |
| 97  | 16.415 | 2482 | 2487 | 2496 | rVB  | 101377 | 198834 | 9.22%  | 0.639% |
| 98  | 16.621 | 2517 | 2522 | 2533 | rVB3 | 42037  | 126310 | 5.86%  | 0.406% |
| 99  | 18.198 | 2784 | 2790 | 2792 | rBV3 | 21841  | 39273  | 1.82%  | 0.126% |
| 100 | 19.133 | 2942 | 2949 | 2955 | rBV7 | 10885  | 32938  | 1.53%  | 0.106% |

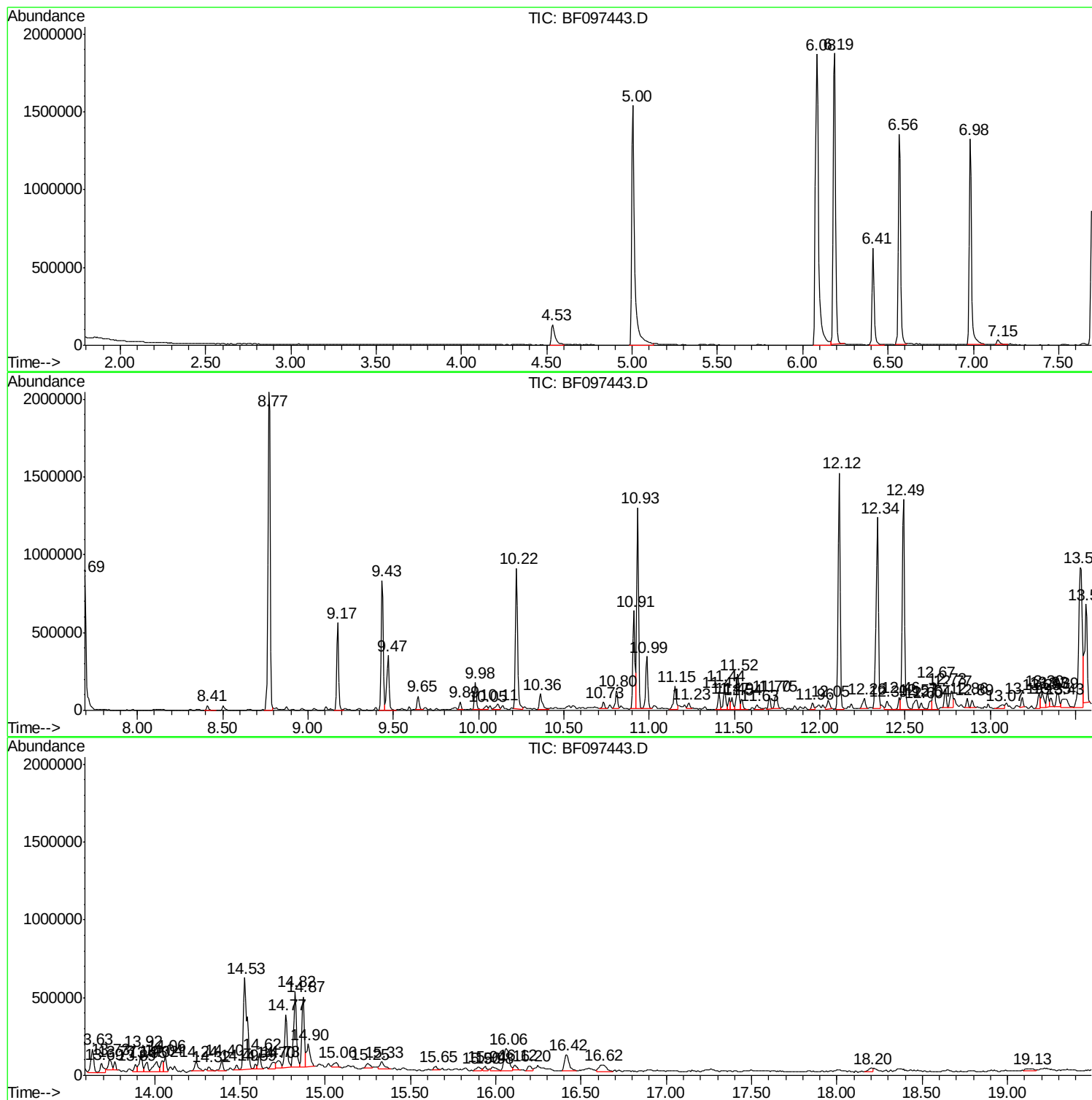
Sum of corrected areas: 31107311

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Instrument :  
BNA\_F  
ClientSampled :  
TEC-WN

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

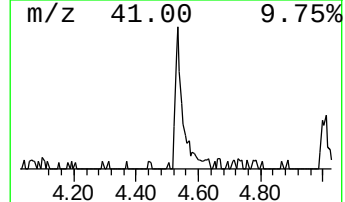
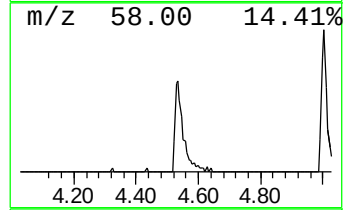
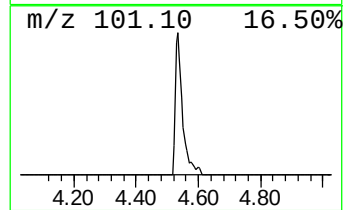
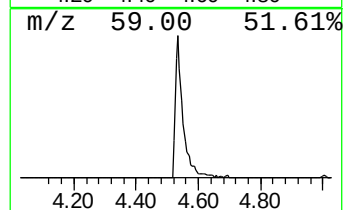
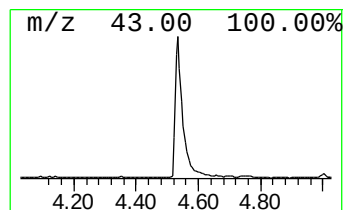
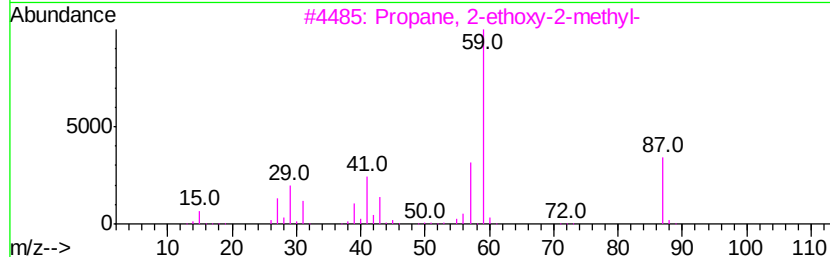
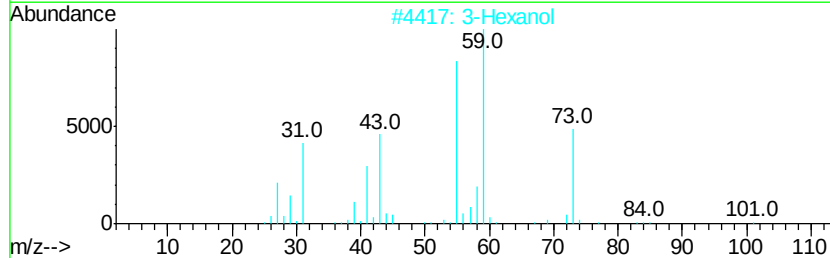
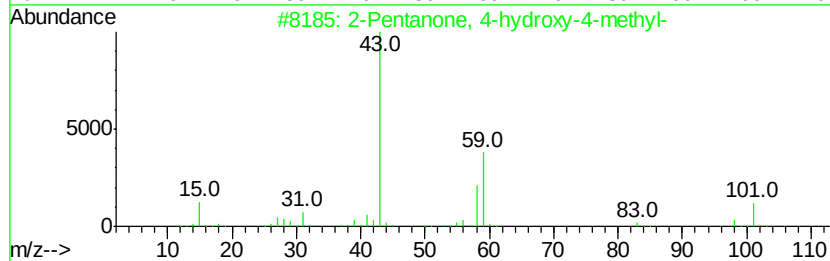
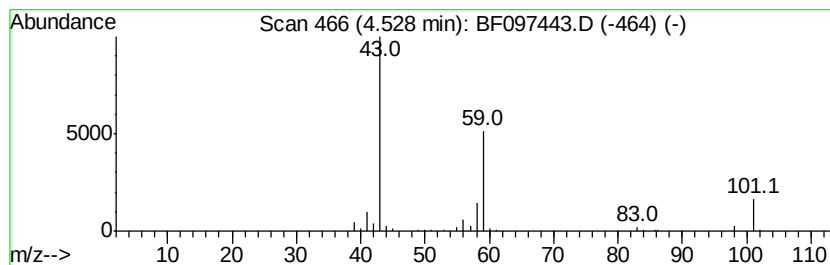
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.53 | 7.41 ng | 211365 | 1,4-Dichlorobenzene-d4 | 6.41 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2 | 50   |
| 2    |    |   | 3-Hexanol                           | 102 | C6H14O  | 000623-37-0 | 33   |
| 3    |    |   | Propane, 2-ethoxy-2-methyl-         | 102 | C6H14O  | 000637-92-3 | 32   |
| 4    |    |   | 2-Hexanol, 2-methyl-                | 116 | C7H16O  | 000625-23-0 | 28   |
| 5    |    |   | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2 | 000540-88-5 | 28   |



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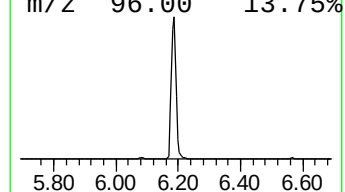
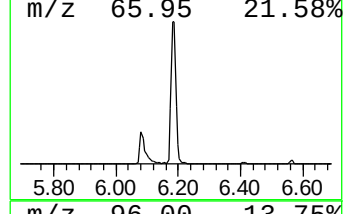
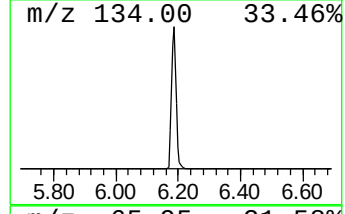
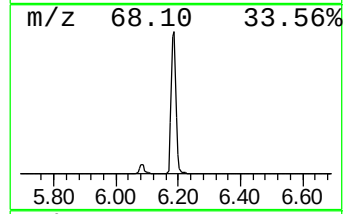
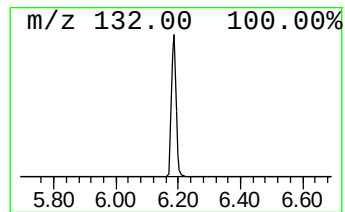
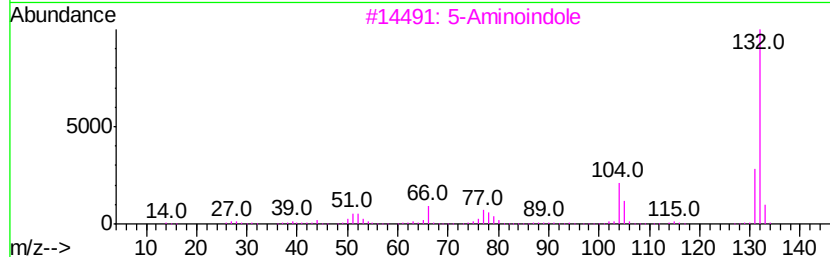
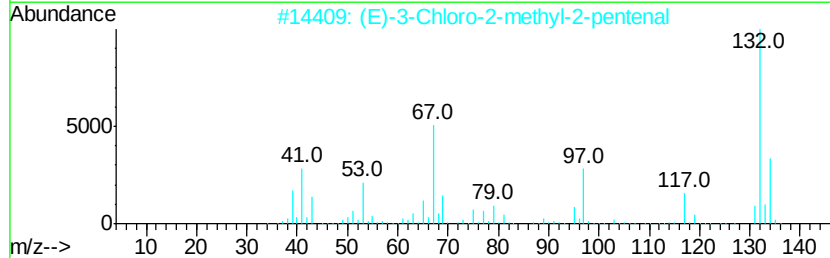
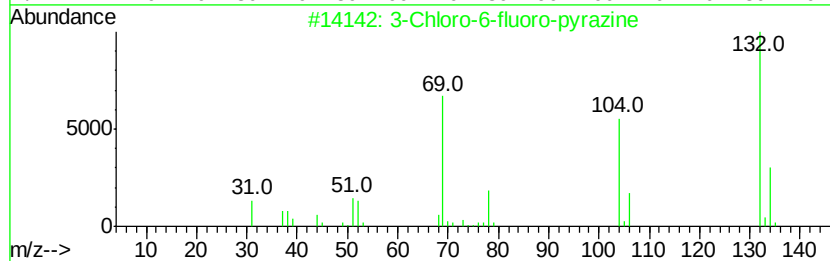
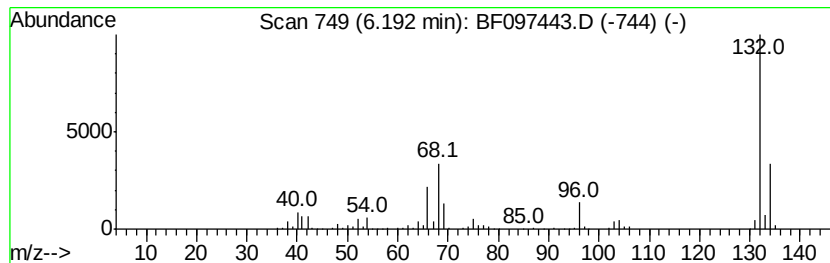
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown6.19 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.19 | 68.13 ng | 1942850 | 1,4-Dichlorobenzene-d4 | 6.41 |

| 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    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 5    |    | 1,3,5-Trifluorobenzene           | 132 | C6H3F3    | 000372-38-3  | 9    |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
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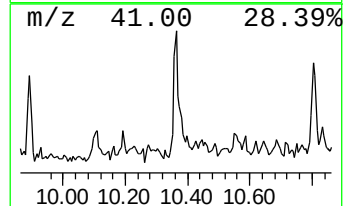
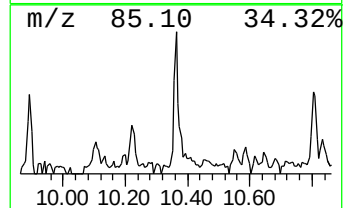
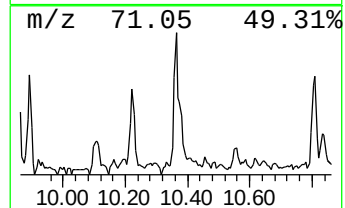
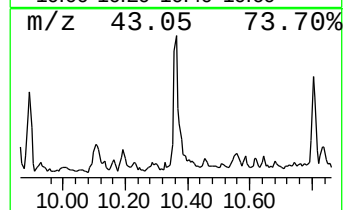
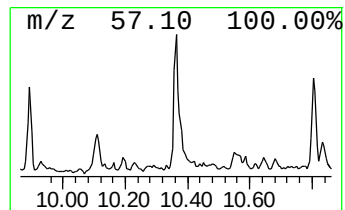
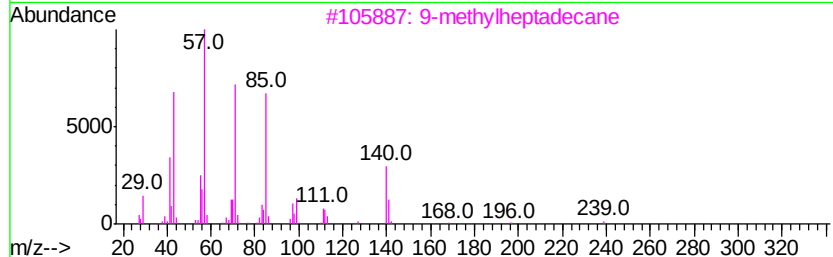
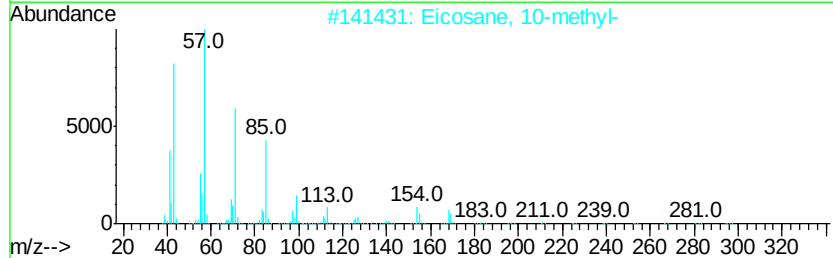
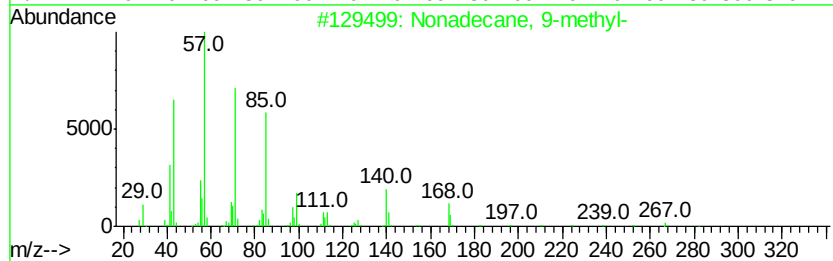
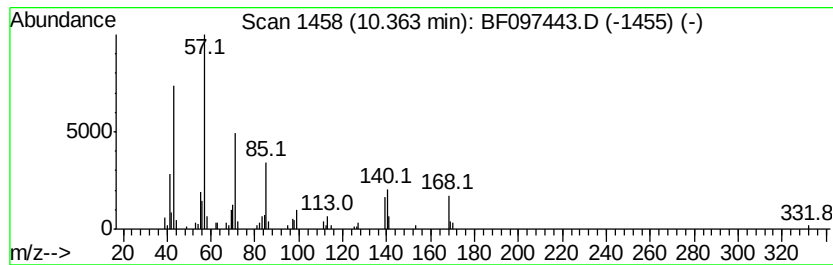
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ClientSampleId :  
TEC-WN

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Nonadecane, 9-methyl- Concentration Rank 9

| R.T.      | EstConc                      | Area   | Relative to ISTD | R.T.        |      |
|-----------|------------------------------|--------|------------------|-------------|------|
| 10.36     | 4.02 ng                      | 120363 | Phenanthrene-d10 | 10.91       |      |
| Hit# of 5 | Tentative ID                 | MW     | MolForm          | CAS#        | Qual |
| 1         | Nonadecane, 9-methyl-        | 282    | C20H42           | 013287-24-6 | 80   |
| 2         | Eicosane, 10-methyl-         | 296    | C21H44           | 054833-23-7 | 62   |
| 3         | 9-methylheptadecane          | 254    | C18H38           | 026741-18-4 | 58   |
| 4         | Heptacosane                  | 380    | C27H56           | 000593-49-7 | 52   |
| 5         | Dodecane, 2-methyl-6-propyl- | 226    | C16H34           | 055045-08-4 | 52   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
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Operator : SJ/JU  
Sample : I4623-06  
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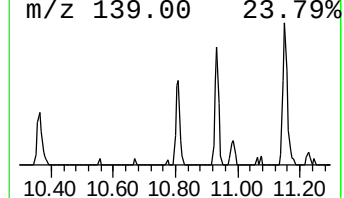
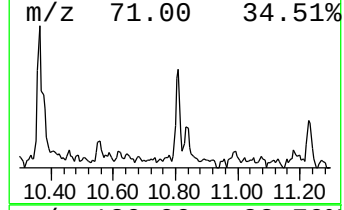
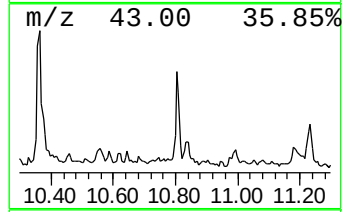
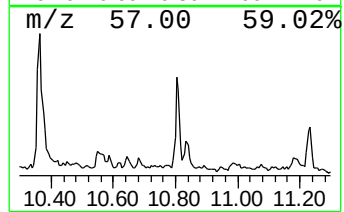
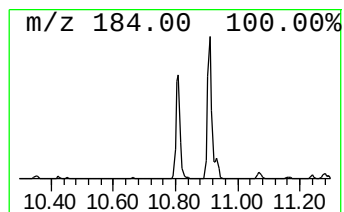
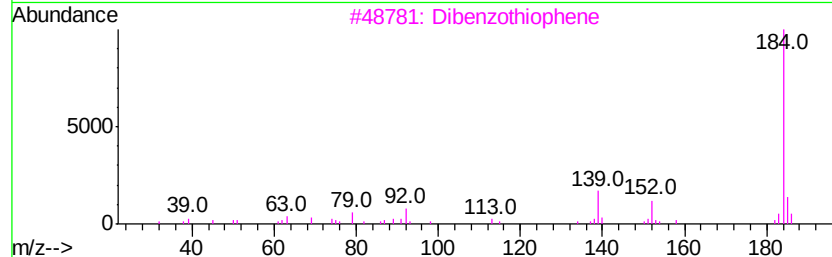
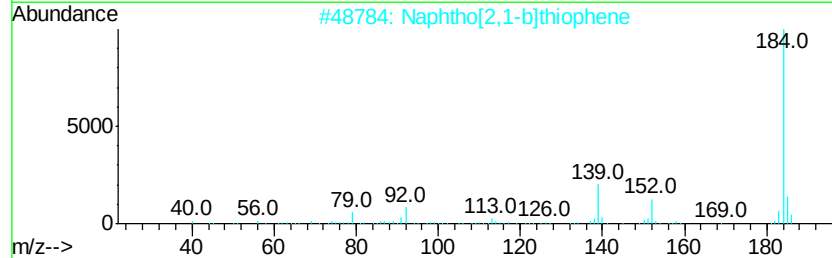
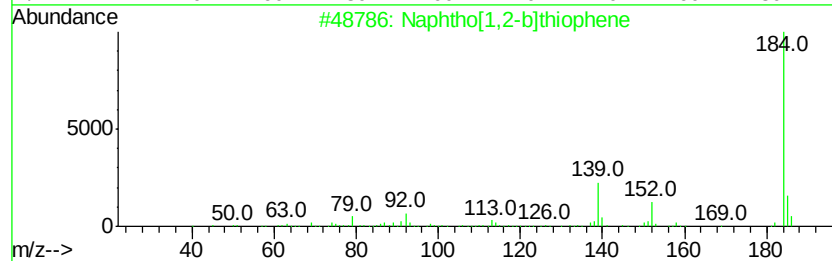
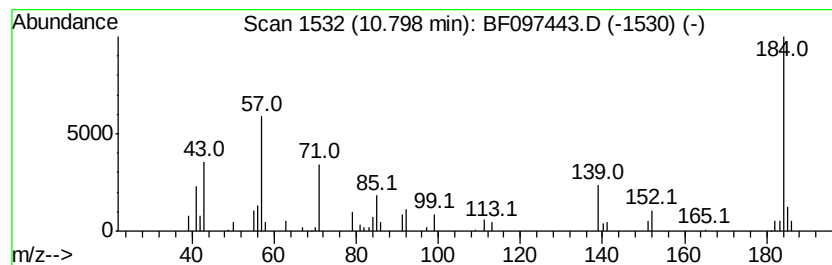
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Naphtho[1,2-b]thiophene Concentration Rank 11

| R.T.      | EstConc                 | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------|--------|------------------|-------------|------|
| 10.80     | 3.35 ng                 | 100335 | Phenanthrene-d10 | 10.91       |      |
| Hit# of 5 | Tentative ID            | MW     | MolForm          | CAS#        | Qual |
| 1         | Naphtho[1,2-b]thiophene | 184    | C12H8S           | 000234-41-3 | 95   |
| 2         | Naphtho[2,1-b]thiophene | 184    | C12H8S           | 000233-02-3 | 95   |
| 3         | Dibenzothiophene        | 184    | C12H8S           | 000132-65-0 | 95   |
| 4         | Naphtho[2,3-b]thiophene | 184    | C12H8S           | 000268-77-9 | 94   |
| 5         | Azuleno(2,1-b)thiophene | 184    | C12H8S           | 000248-13-5 | 90   |





Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

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

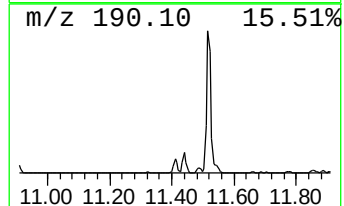
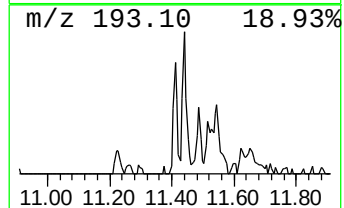
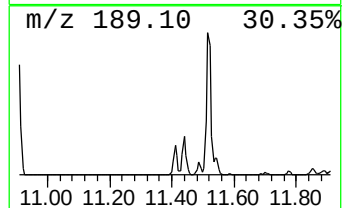
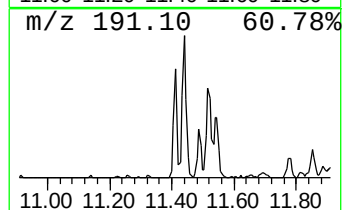
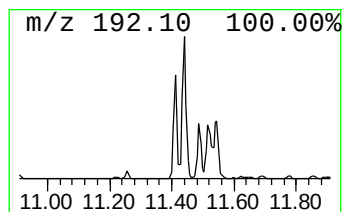
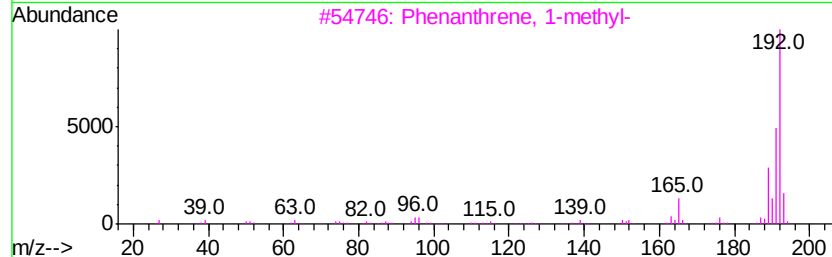
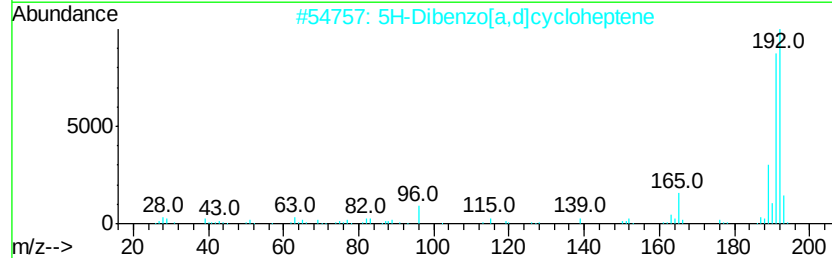
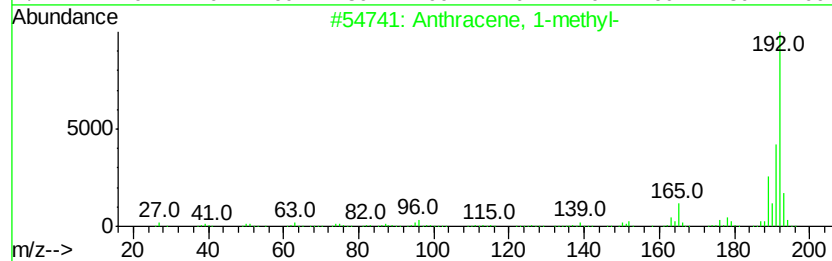
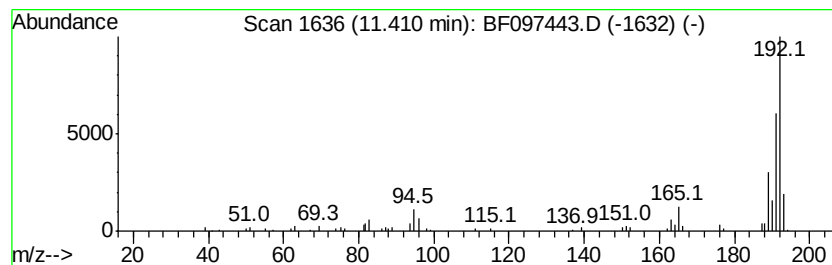
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Anthracene, 1-methyl- Concentration Rank 15

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.41 | 3.00 ng | 89744 | Phenanthrene-d10 | 10.91 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 1-methyl-       | 192 | C15H12  | 000610-48-0 | 95   |
| 2    |    | 5H-Dibenzo[a,d]cycloheptene | 192 | C15H12  | 000256-81-5 | 93   |
| 3    |    | Phenanthrene, 1-methyl-     | 192 | C15H12  | 000832-69-9 | 93   |
| 4    |    | Phenanthrene, 2-methyl-     | 192 | C15H12  | 002531-84-2 | 93   |
| 5    |    | Anthracene, 2-methyl-       | 192 | C15H12  | 000613-12-7 | 91   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

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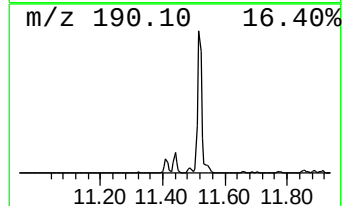
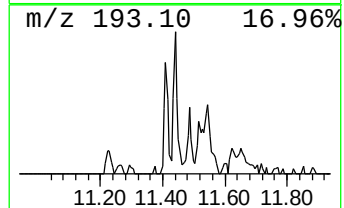
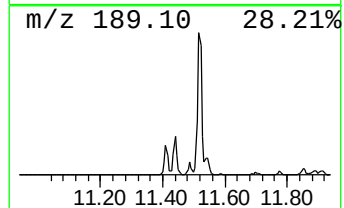
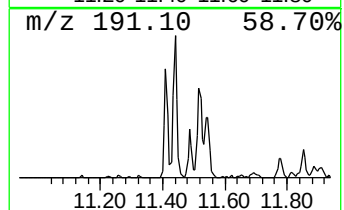
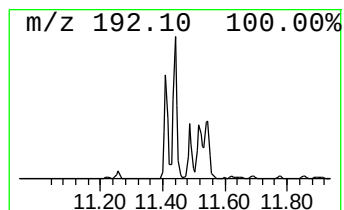
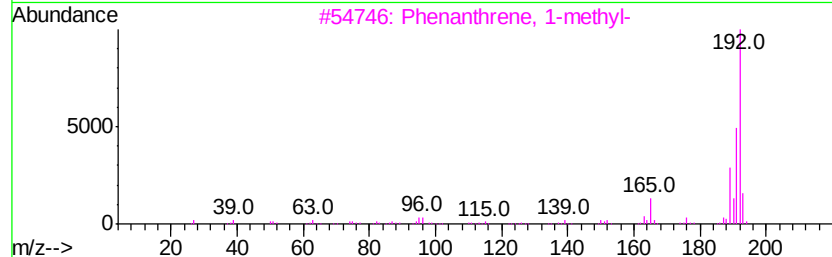
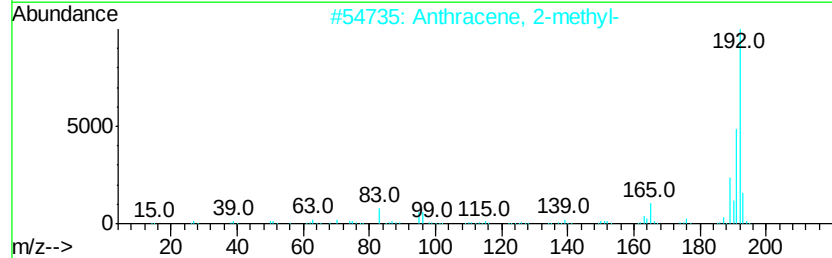
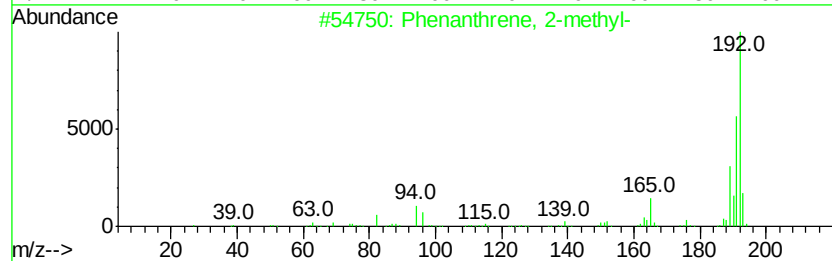
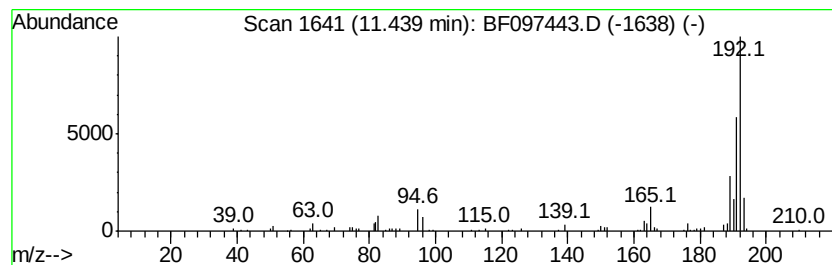
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.44 | 4.76 ng | 142425 | Phenanthrene-d10 | 10.91 |

| Hit# of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------|-----|---------|-------------|------|
| 1       |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 98   |
| 2       |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 96   |
| 3       |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 96   |
| 4       |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 96   |
| 5       |   | Phenanthrene, 4-methyl- | 192 | C15H12  | 000832-64-4 | 93   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
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Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

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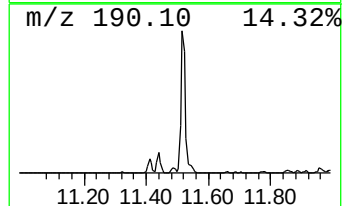
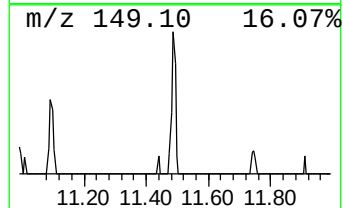
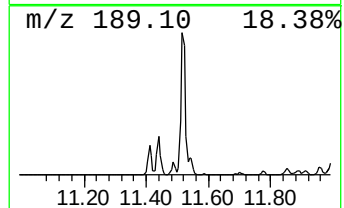
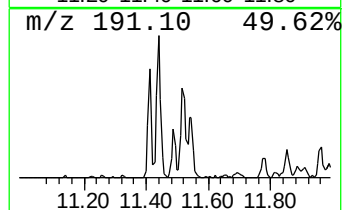
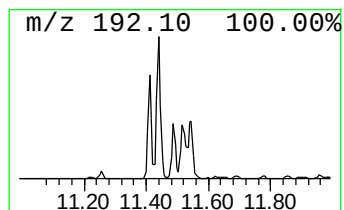
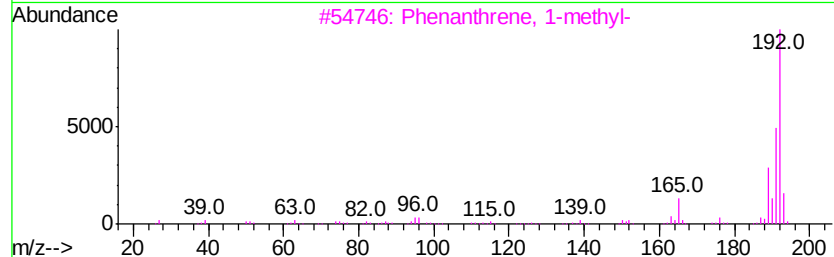
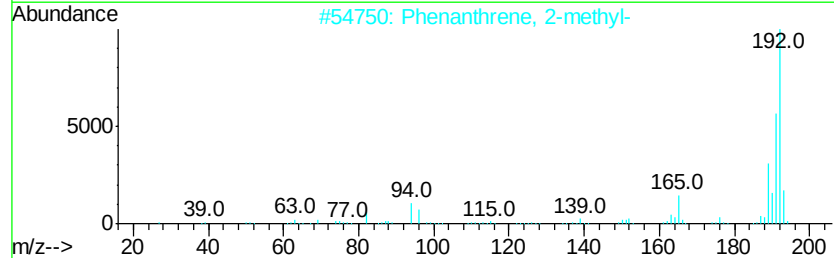
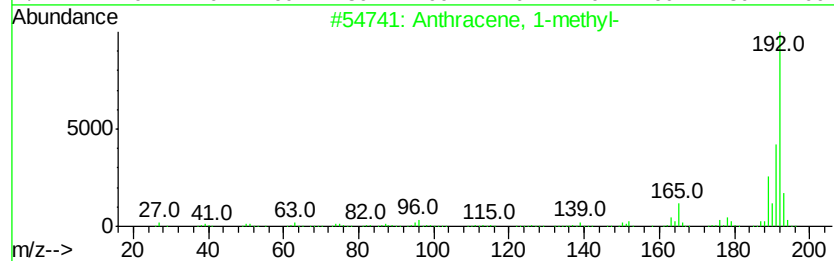
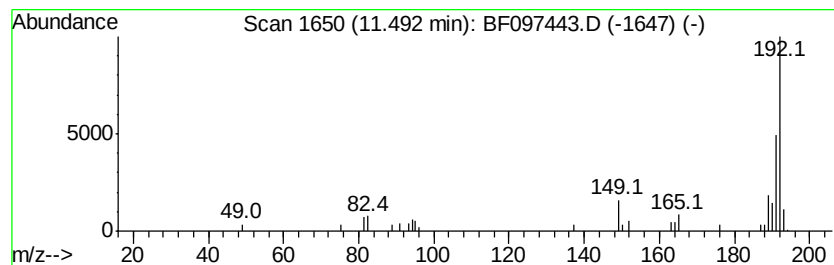
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 19

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.49 | 2.40 ng | 71863 | Phenanthrene-d10 | 10.91 |

| Hit# | of | Tentative ID                           | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 1-methyl-                  | 192 | C15H12  | 000610-48-0 | 97   |
| 2    |    | Phenanthrene, 2-methyl-                | 192 | C15H12  | 002531-84-2 | 96   |
| 3    |    | Phenanthrene, 1-methyl-                | 192 | C15H12  | 000832-69-9 | 96   |
| 4    |    | 1H-Cyclopropa[1,1]phenanthrene, 1a,... | 192 | C15H12  | 000949-41-7 | 95   |
| 5    |    | Anthracene, 2-methyl-                  | 192 | C15H12  | 000613-12-7 | 94   |



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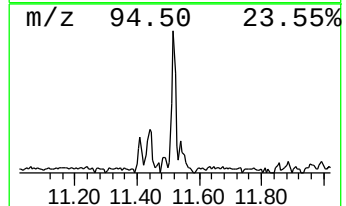
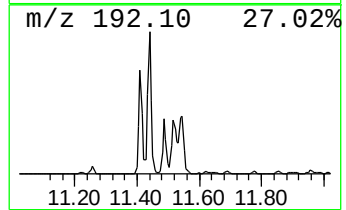
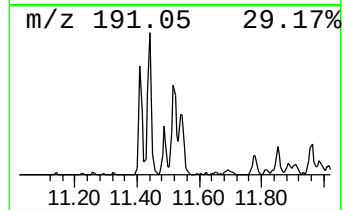
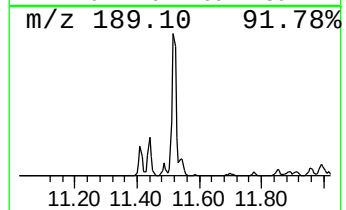
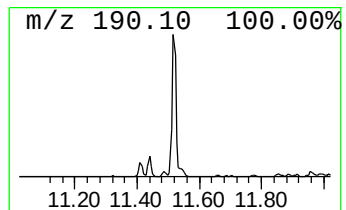
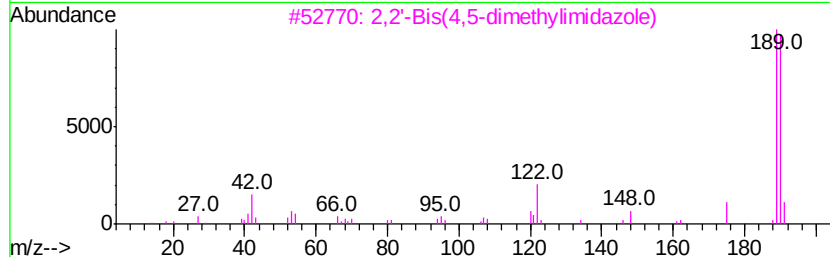
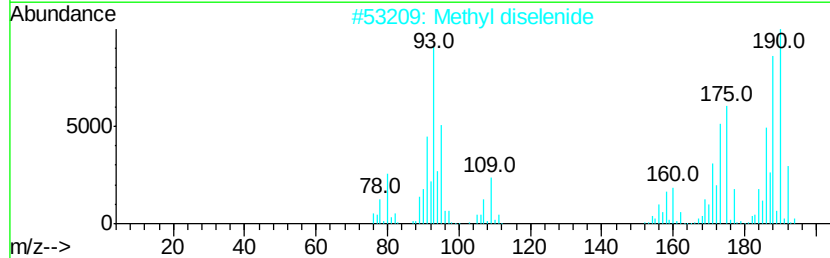
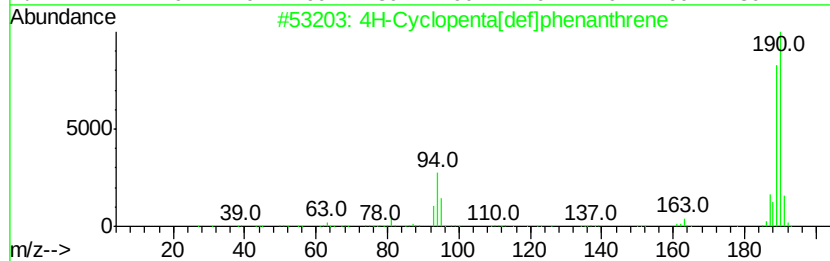
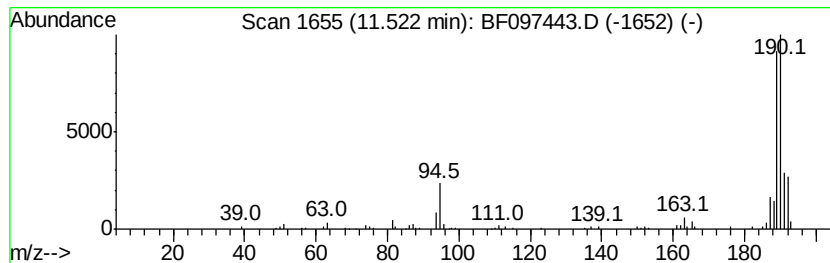
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 5

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.52     | 7.31 ng                             | 218762 | Phenanthrene-d10 | 10.91       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | 4H-Cyclopenta[def]phenanthrene      | 190    | C15H10           | 000203-64-5 | 93   |
| 2         | Methyl diselenide                   | 190    | C2H6Se2          | 007101-31-7 | 41   |
| 3         | 2,2'-Bis(4,5-dimethylimidazole)     | 190    | C10H14N4         | 069286-06-2 | 38   |
| 4         | 6,7-Methylenedioxy-4(3H)-quinazo... | 190    | C9H6N2O3         | 028310-11-4 | 28   |
| 5         | 1,4-Naphthalenedione, 2,5-dihydr... | 190    | C10H6O4          | 004923-55-1 | 25   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

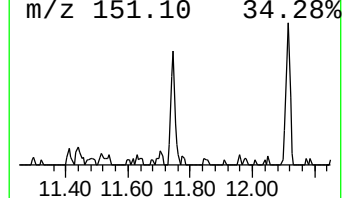
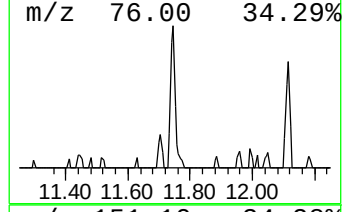
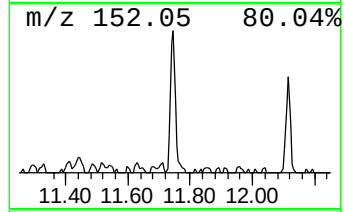
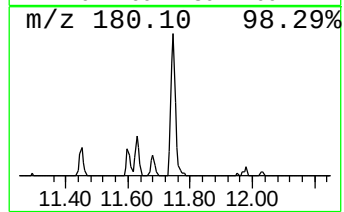
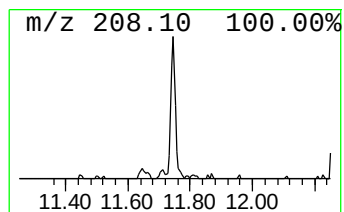
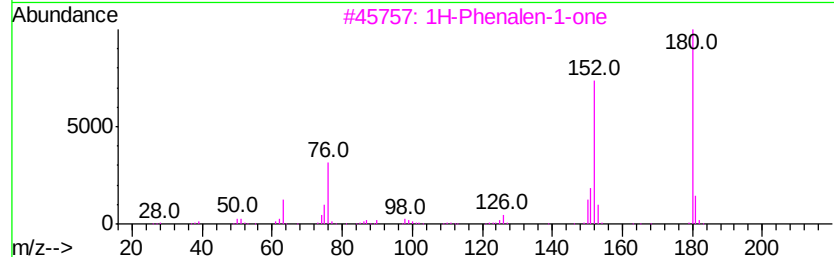
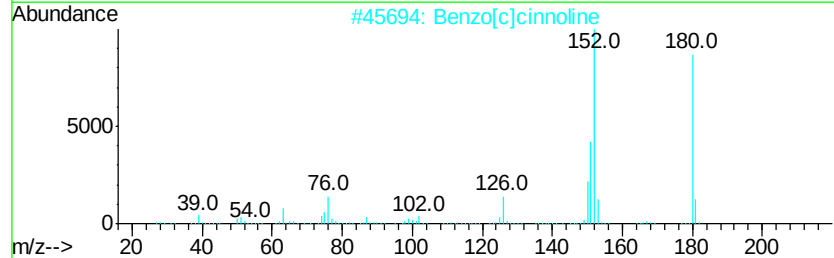
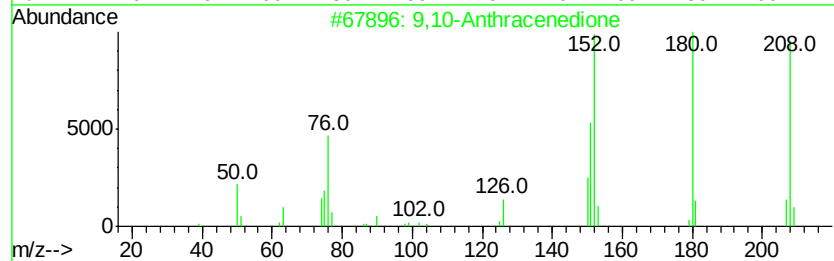
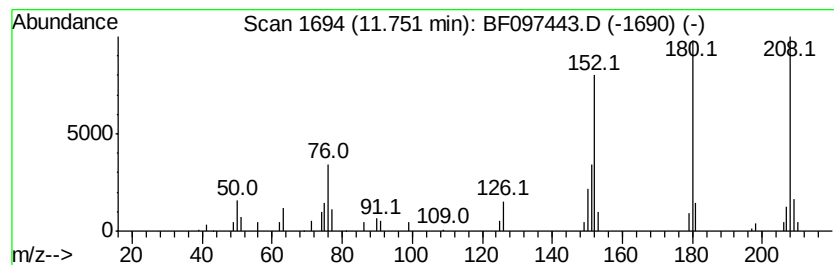
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 9,10-Anthracenedione Concentration Rank 17

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 11.75 | 2.70 ng | 80749 | Phenanthrene-d10 | 10.91 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 97   |
| 2    |    | Benzo[c]cinnoline    | 180 | C12H8N2 | 000230-17-1 | 92   |
| 3    |    | 1H-Phenalen-1-one    | 180 | C13H8O  | 000548-39-0 | 78   |
| 4    |    | 9H-Fluoren-9-one     | 180 | C13H8O  | 000486-25-9 | 62   |
| 5    |    | 1,4-Anthracenedione  | 208 | C14H8O2 | 000635-12-1 | 49   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

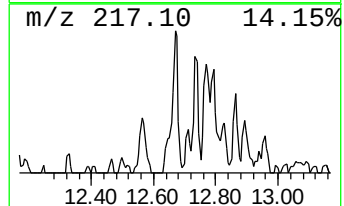
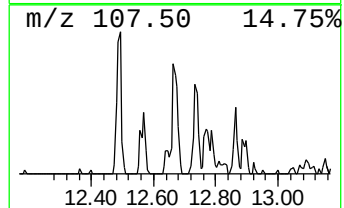
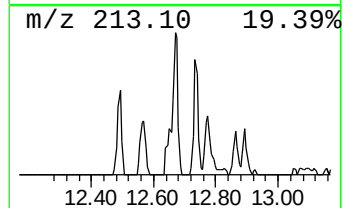
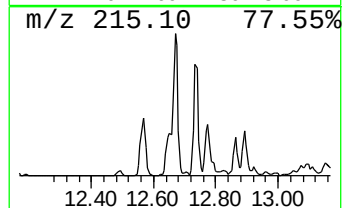
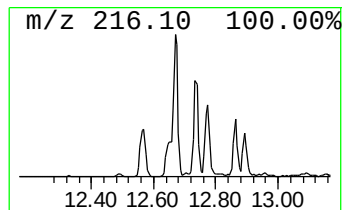
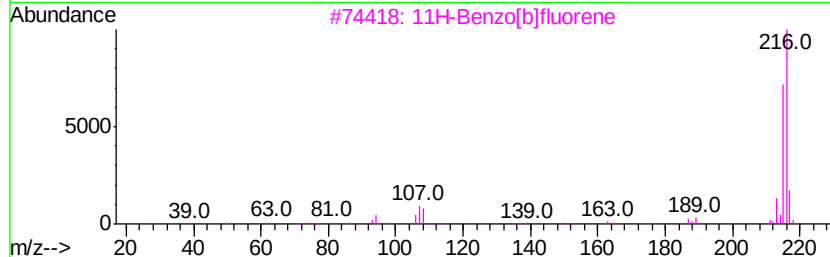
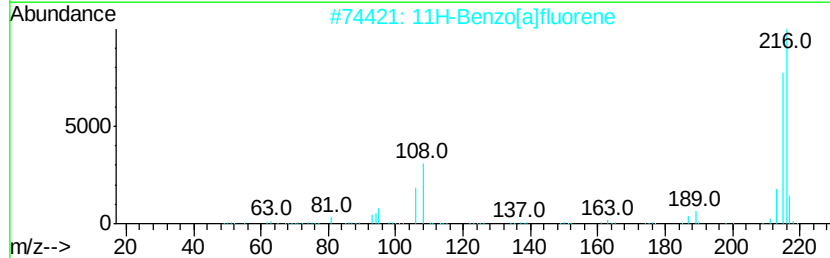
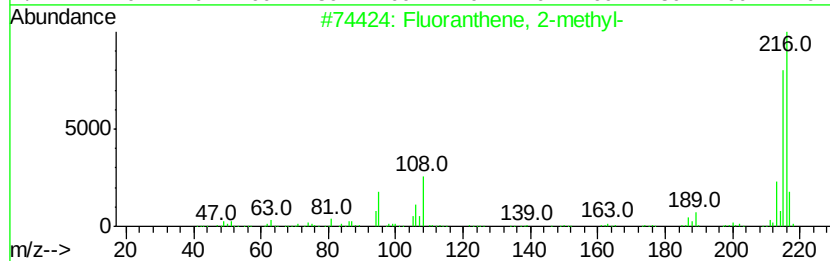
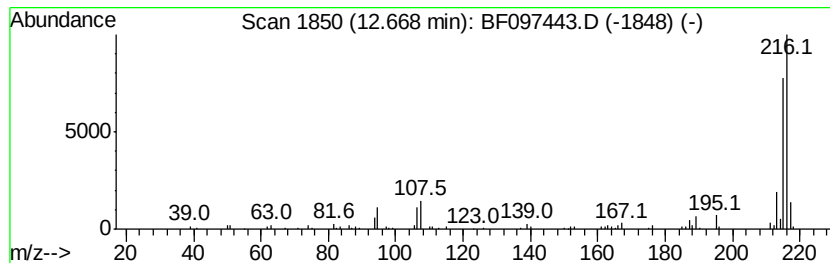
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 Fluoranthene, 2-methyl- Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.67 | 2.38 ng | 167749 | Chrysene-d12     | 13.53 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 95   |
| 2    |    | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 93   |
| 3    |    | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 90   |
| 4    |    | Pvrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 87   |
| 5    |    | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 81   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

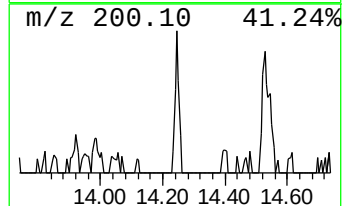
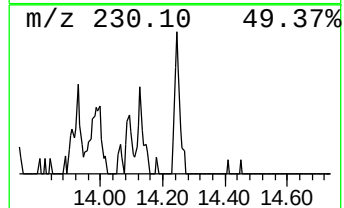
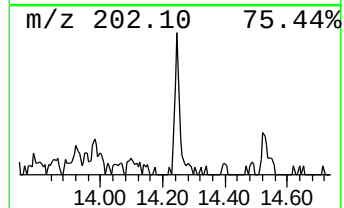
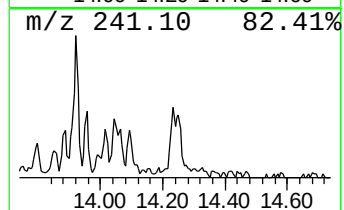
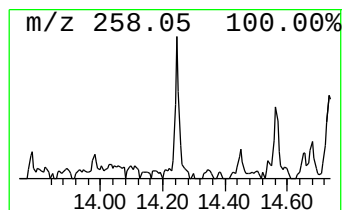
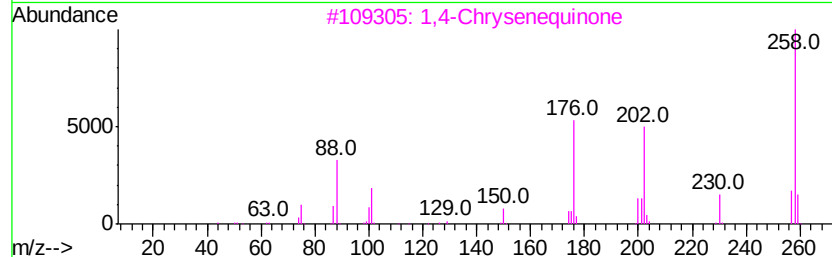
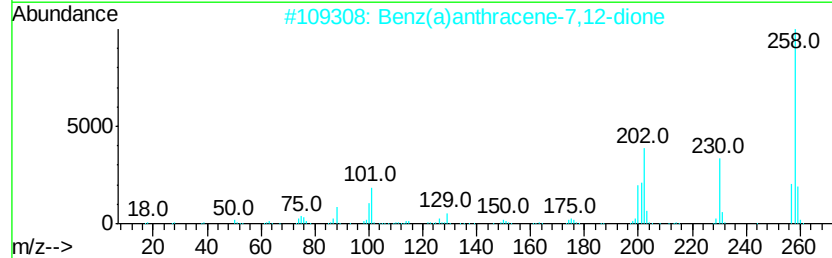
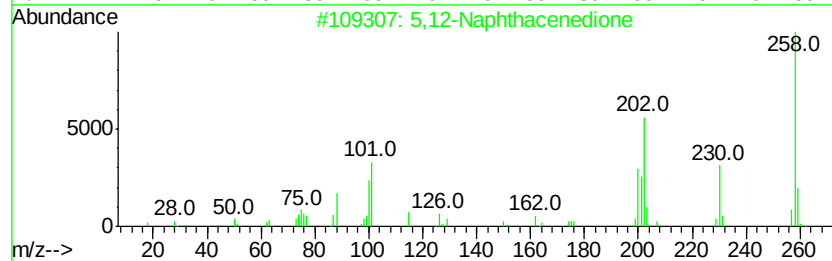
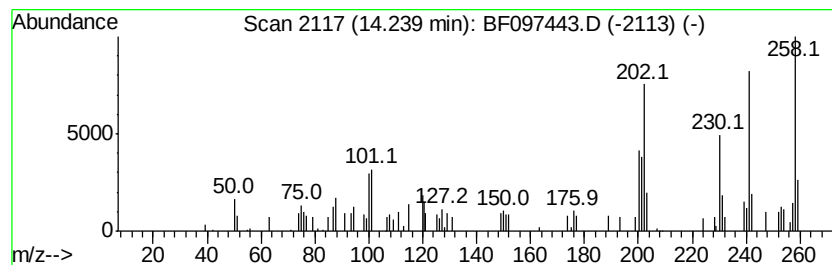
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 5,12-Naphthacenedione Concentration Rank 10

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.24 | 3.68 ng | 90246 | Perylene-d12     | 14.87 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 5,12-Naphthacenedione               | 258 | C18H10O2 | 001090-13-7 | 95   |
| 2    |    |   | Benz(a)anthracene-7,12-dione        | 258 | C18H10O2 | 002498-66-0 | 90   |
| 3    |    |   | 1,4-Chrysenequinone                 | 258 | C18H10O2 | 100900-16-1 | 50   |
| 4    |    |   | Benzo[1,2-b:4,3-b']bisbenzofuran    | 258 | C18H10O2 | 000192-93-8 | 49   |
| 5    |    |   | 4,5,7,8,9,10-Hexahydrobenzo[a]py... | 258 | C20H18   | 073712-75-1 | 43   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

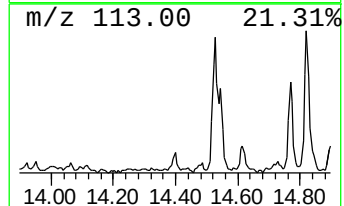
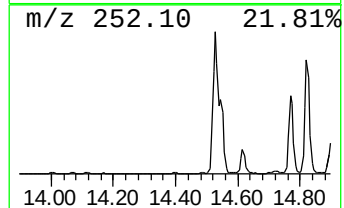
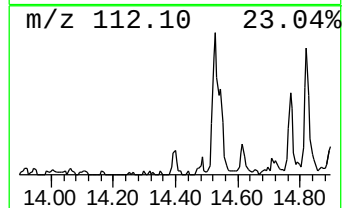
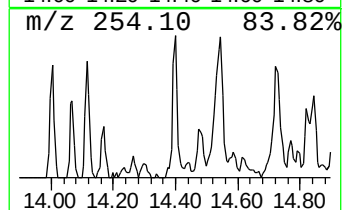
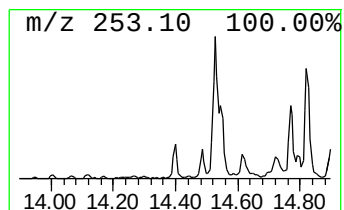
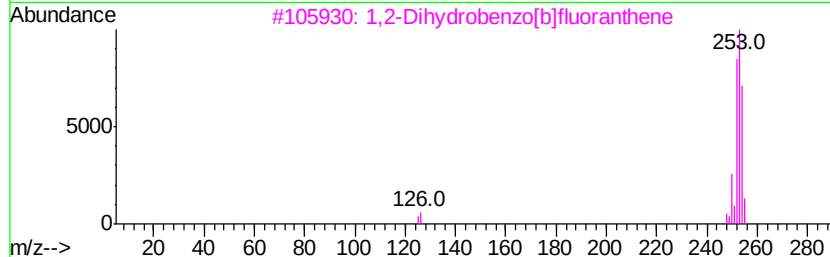
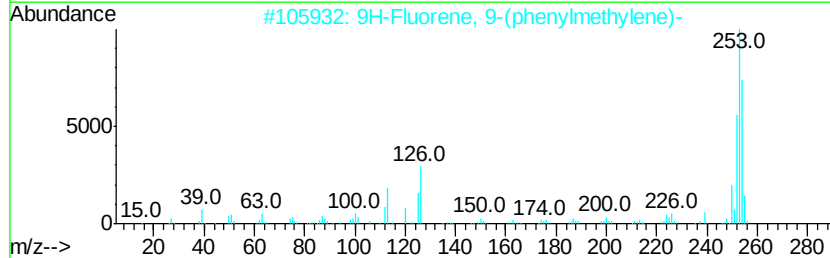
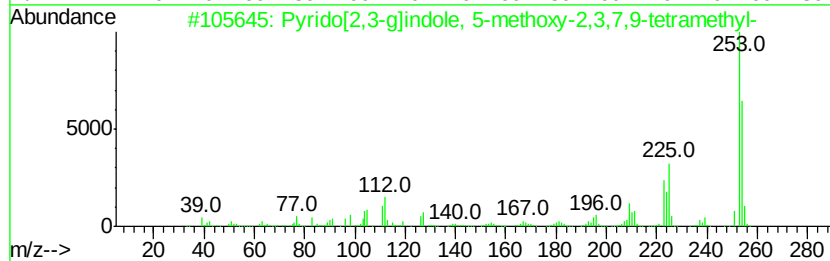
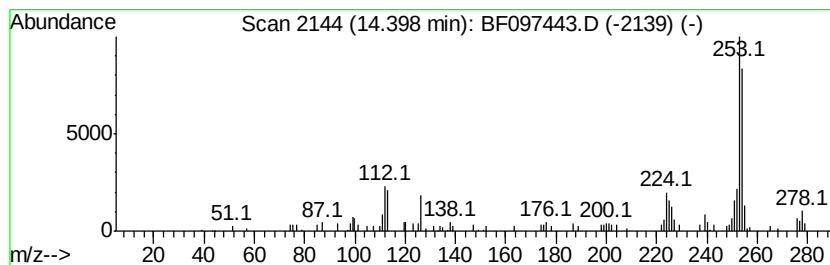
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TIC Integration Parameters: LSCINT.P

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Peak Number 13 Pyrido[2,3-g]indole, 5-meth... Concentration Rank 16

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.40 | 2.96 ng | 72543 | Perylene-d12     | 14.87 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Pyrido[2,3-g]indole, 5-methoxy-2... | 254 | C16H18N2O | 201991-45-9 | 64   |
| 2    |    | 9H-Fluorene, 9-(phenylmethylene)-   | 254 | C20H14    | 001836-87-9 | 62   |
| 3    |    | 1,2-Dihydrobenzo[b]fluoranthene     | 254 | C20H14    | 100516-13-0 | 58   |
| 4    |    | 4,6'-BiazulenyI                     | 254 | C20H14    | 094154-49-1 | 53   |
| 5    |    | Benz(a)anthracene-7-carbonitrile    | 253 | C19H11N   | 007476-08-6 | 50   |





Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

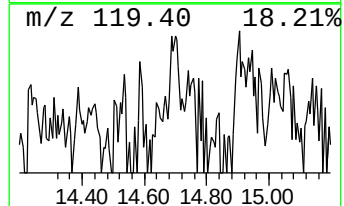
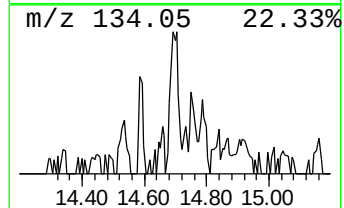
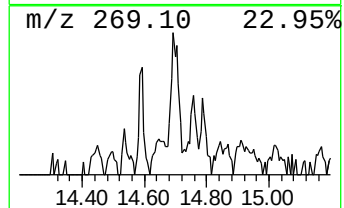
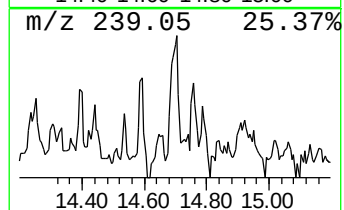
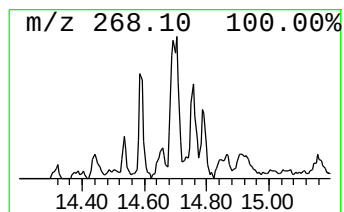
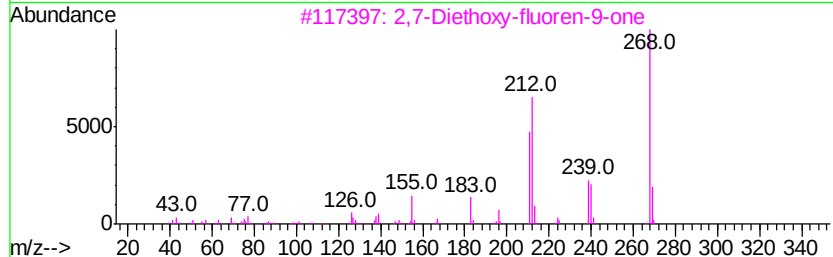
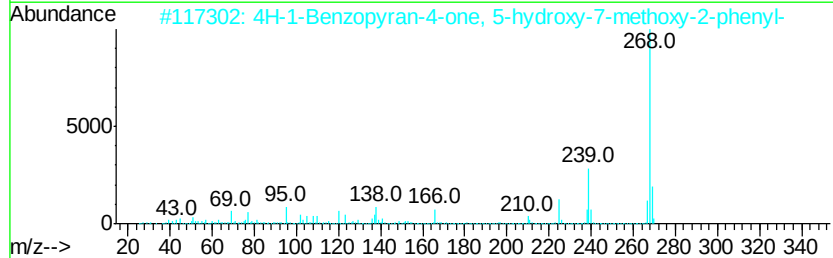
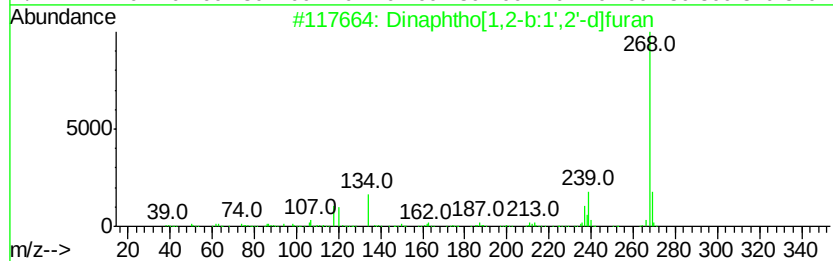
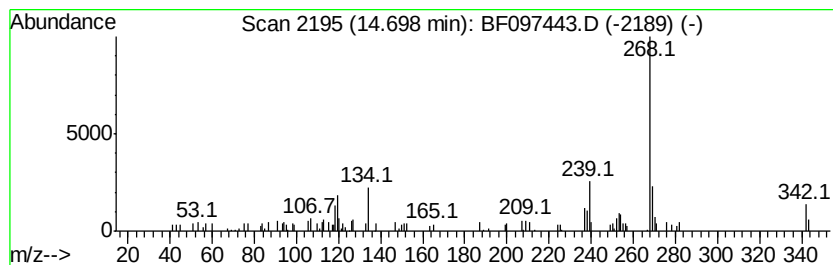
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Dinaphtho[1,2-b:1',2'-d]furan Concentration Rank 12

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.70 | 3.29 ng | 80672 | Perylene-d12     | 14.87 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Dinaphtho[1,2-b:1',2'-d]furan       | 268 | C20H12O   | 000207-93-2  | 87   |
| 2    |    | 4H-1-Benzopyran-4-one, 5-hydroxy... | 268 | C16H12O4  | 000520-28-5  | 53   |
| 3    |    | 2,7-Diethoxy-fluoren-9-one          | 268 | C17H16O3  | 1000317-69-6 | 52   |
| 4    |    | 7,9-Diamino-5,6-dihydronaphtho[2... | 268 | C14H12N4S | 037071-29-7  | 52   |
| 5    |    | Naphthalene, 1,1'-methylenebis-     | 268 | C21H16    | 000607-50-1  | 47   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

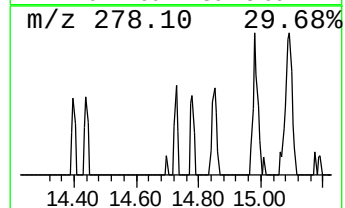
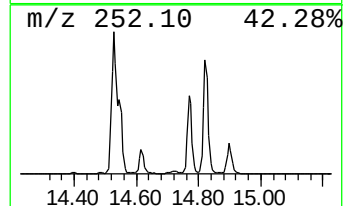
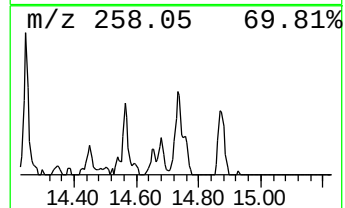
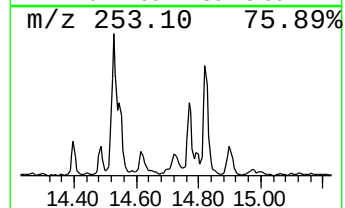
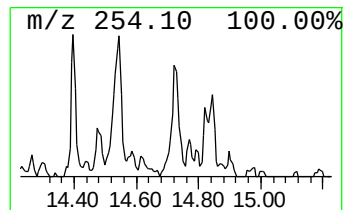
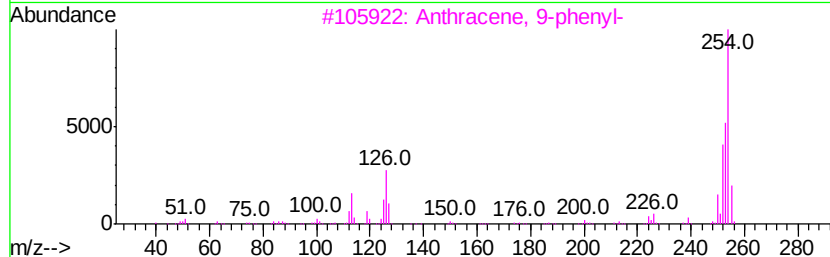
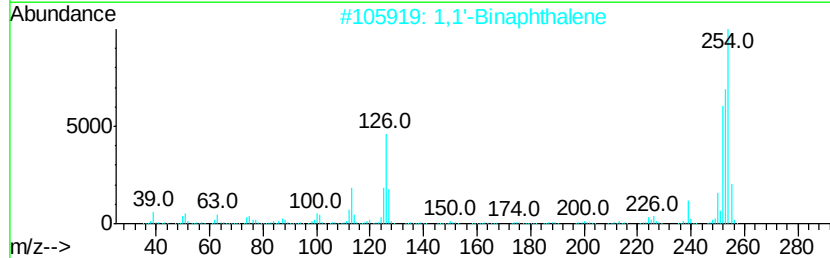
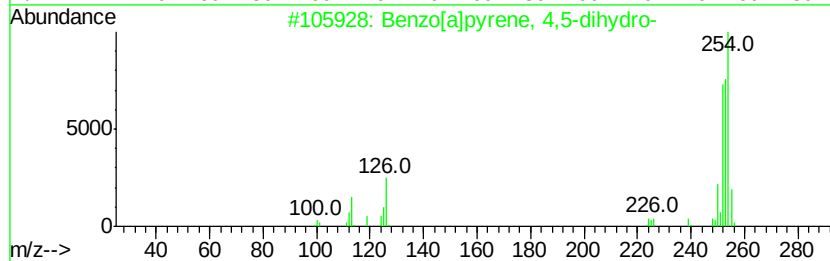
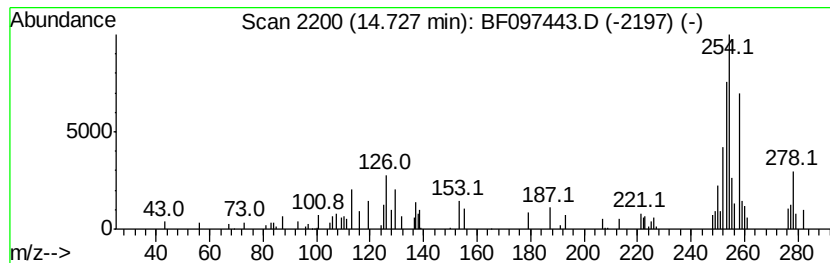
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzo[a]pyrene, 4,5-dihydro- Concentration Rank 13

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.73 | 3.08 ng | 75492 | Perylene-d12     | 14.87 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[a]pyrene, 4,5-dihydro- | 254 | C20H14  | 057652-66-1 | 64   |
| 2    |    | 1,1'-Binaphthalene           | 254 | C20H14  | 000604-53-5 | 50   |
| 3    |    | Anthracene, 9-phenyl-        | 254 | C20H14  | 000602-55-1 | 50   |
| 4    |    | Benzo(a)pyrene, 7,8-dihydro- | 254 | C20H14  | 017573-23-8 | 50   |
| 5    |    | 1,2'-Binaphthalene           | 254 | C20H14  | 004325-74-0 | 45   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

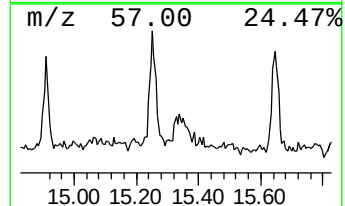
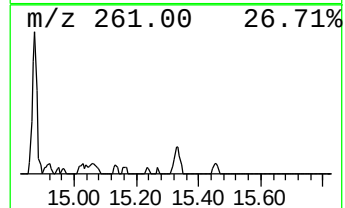
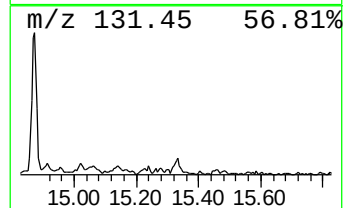
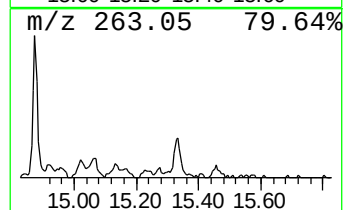
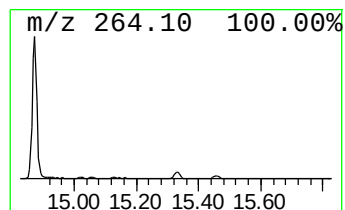
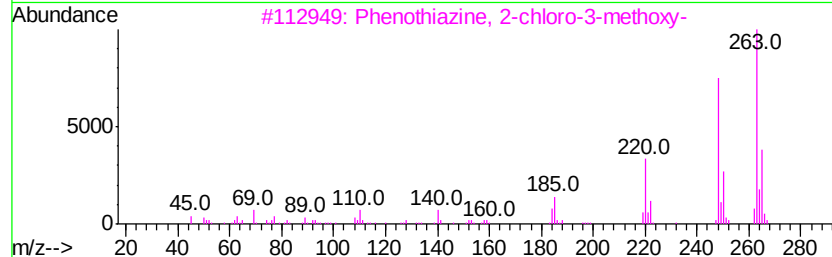
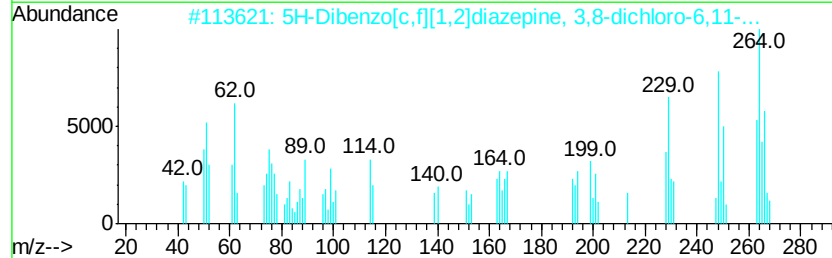
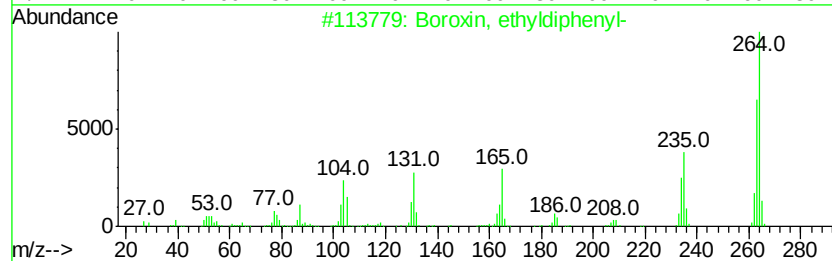
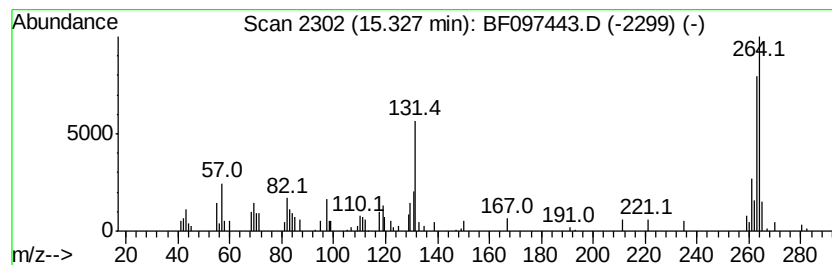
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 unknown15.33 Concentration Rank 14

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 15.33 | 3.06 ng | 75043 | Perylene-d12     | 14.87 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Boroxin, ethyldiphenyl-             | 264 | C14H15B3O3  | 1000151-82-0 | 40   |
| 2    |    | 5H-Dibenzo[c,f][1,2]diazepine, 3... | 264 | C13H10Cl2N2 | 000955-66-8  | 27   |
| 3    |    | Phenothiazine, 2-chloro-3-methoxy-  | 263 | C13H10ClNO3 | 017800-08-7  | 27   |
| 4    |    | 1,3-Butanedione, 1-phenyl-2-(ph...  | 265 | C17H15NO2   | 101441-99-0  | 22   |
| 5    |    | 2-Oxo-6-phenyl-4-(4-hydroxypheny... | 264 | C16H12N2O2  | 161200-20-0  | 16   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF080717\  
Data File : BF097443.D  
Acq On : 7 Aug 2017 21:48  
Operator : SJ/JU  
Sample : I4623-06  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TEC-WN

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

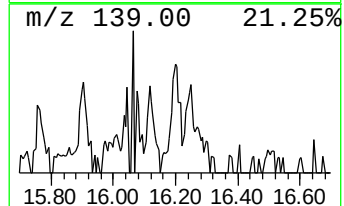
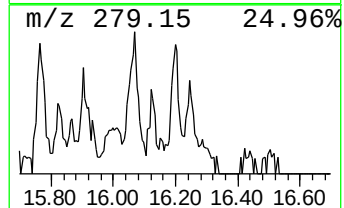
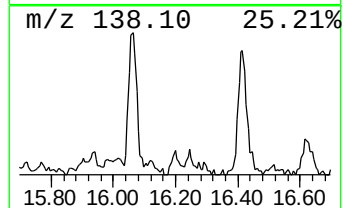
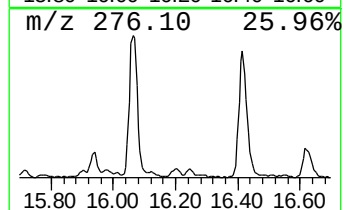
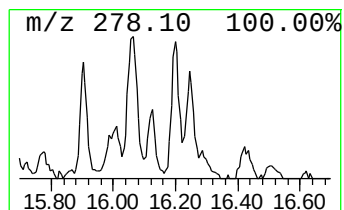
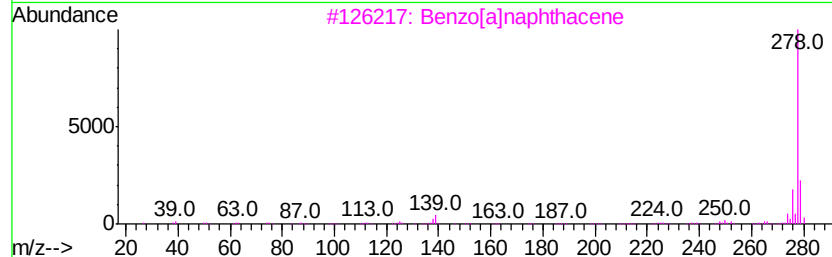
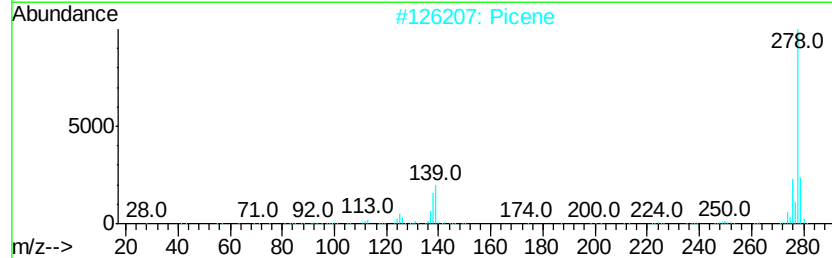
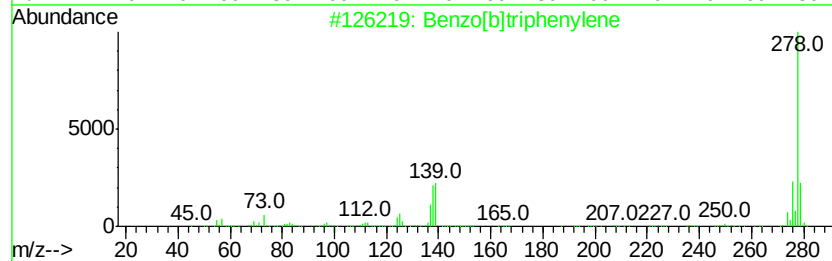
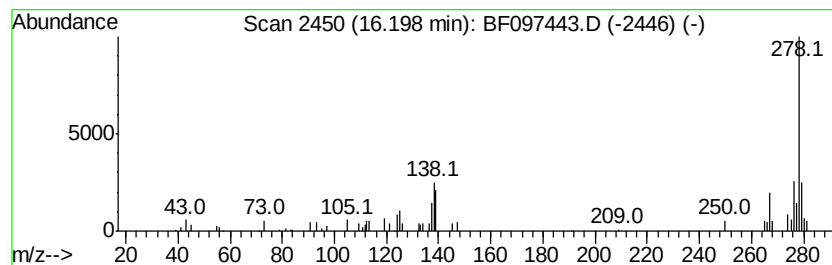
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Benzo[b]triphenylene Concentration Rank 18

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.20 | 2.48 ng | 60808 | Perylene-d12     | 14.87 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[b]triphenylene  | 278 | C22H14  | 000215-58-7 | 96   |
| 2    |    | Picene                | 278 | C22H14  | 000213-46-7 | 68   |
| 3    |    | Benzo[a]naphthacene   | 278 | C22H14  | 000226-88-0 | 58   |
| 4    |    | Dibenz[a,h]anthracene | 278 | C22H14  | 000053-70-3 | 52   |
| 5    |    | Pentacene             | 278 | C22H14  | 000135-48-8 | 52   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF080717\  
 Data File : BF097443.D  
 Acq On : 7 Aug 2017 21:48  
 Operator : SJ/JU  
 Sample : I4623-06  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TEC-WN

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF072517.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... | 4.53  | 7.4     | ng    | 211365   | 1                     | 6.41  | 570308  | 20.0 |
| unknown6.19          | 6.19  | 68.1    | ng    | 1942850  | 1                     | 6.41  | 570308  | 20.0 |
| Nonadecane, 9-met... | 10.36 | 4.0     | ng    | 120363   | 4                     | 10.91 | 598652  | 20.0 |
| Naphtho[1,2-b]thi... | 10.80 | 3.4     | ng    | 100335   | 4                     | 10.91 | 598652  | 20.0 |
| Anthracene, 1-met... | 11.41 | 3.0     | ng    | 89744    | 4                     | 10.91 | 598652  | 20.0 |
| Phenanthrene, 2-m... | 11.44 | 4.8     | ng    | 142425   | 4                     | 10.91 | 598652  | 20.0 |
| Phenanthrene, 1-m... | 11.49 | 2.4     | ng    | 71863    | 4                     | 10.91 | 598652  | 20.0 |
| 4H-Cyclopenta[def... | 11.52 | 7.3     | ng    | 218762   | 4                     | 10.91 | 598652  | 20.0 |
| 9,10-Anthracenedione | 11.75 | 2.7     | ng    | 80749    | 4                     | 10.91 | 598652  | 20.0 |
| Fluoranthene, 2-m... | 12.67 | 2.4     | ng    | 167749   | 5                     | 13.53 | 1407630 | 20.0 |
| 5,12-Naphthacened... | 14.24 | 3.7     | ng    | 90246    | 6                     | 14.87 | 490858  | 20.0 |
| Pyrido[2,3-g]indo... | 14.40 | 3.0     | ng    | 72543    | 6                     | 14.87 | 490858  | 20.0 |
| Dinaphtho[1,2-b:1... | 14.70 | 3.3     | ng    | 80672    | 6                     | 14.87 | 490858  | 20.0 |
| Benzo[a]pyrene, 4... | 14.73 | 3.1     | ng    | 75492    | 6                     | 14.87 | 490858  | 20.0 |
| unknown15.33         | 15.33 | 3.1     | ng    | 75043    | 6                     | 14.87 | 490858  | 20.0 |
| Benzo[b]triphenylene | 16.20 | 2.5     | ng    | 60808    | 6                     | 14.87 | 490858  | 20.0 |