

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF072920\  
 Data File : BF121118.D  
 Acq On : 29 Jul 2020 23:33  
 Operator : JU/CG  
 Sample : L3436-17 5X  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 17

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

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

Signal : TIC

| peak<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         | 5.416       | 563           | 566         | 581          | rVB      | 542842         | 518625        | 25.20%          | 1.967%        |
| 2         | 6.440       | 737           | 740         | 755          | rBV      | 528934         | 528504        | 25.68%          | 2.005%        |
| 3         | 6.640       | 770           | 774         | 779          | rBV3     | 34663          | 43636         | 2.12%           | 0.166%        |
| 4         | 6.804       | 799           | 802         | 806          | rBV      | 998177         | 853484        | 41.48%          | 3.238%        |
| 5         | 7.069       | 843           | 847         | 853          | rBV2     | 26534          | 37522         | 1.82%           | 0.142%        |
| 6         | 7.363       | 894           | 897         | 902          | rVB      | 333426         | 277442        | 13.48%          | 1.052%        |
| 7         | 7.834       | 974           | 977         | 982          | rBV6     | 16598          | 29488         | 1.43%           | 0.112%        |
| 8         | 8.092       | 1015          | 1021        | 1023         | rBV      | 1207381        | 1081991       | 52.58%          | 4.105%        |
| 9         | 8.110       | 1023          | 1024        | 1027         | rVB      | 205025         | 126708        | 6.16%           | 0.481%        |
| 10        | 8.510       | 1089          | 1092        | 1094         | rBV2     | 26900          | 26014         | 1.26%           | 0.099%        |
| 11        | 8.681       | 1118          | 1121        | 1124         | rBV2     | 37599          | 32220         | 1.57%           | 0.122%        |
| 12        | 8.804       | 1139          | 1142        | 1145         | rVB      | 133258         | 104376        | 5.07%           | 0.396%        |
| 13        | 8.904       | 1155          | 1159        | 1163         | rBV      | 97420          | 100267        | 4.87%           | 0.380%        |
| 14        | 9.110       | 1192          | 1194        | 1199         | rBV3     | 41151          | 36333         | 1.77%           | 0.138%        |
| 15        | 9.163       | 1200          | 1203        | 1208         | rVB      | 843379         | 634305        | 30.83%          | 2.406%        |
| 16        | 9.245       | 1213          | 1217        | 1219         | rBV      | 70884          | 66975         | 3.25%           | 0.254%        |
| 17        | 9.363       | 1234          | 1237        | 1242         | rVB2     | 35295          | 43192         | 2.10%           | 0.164%        |
| 18        | 9.433       | 1242          | 1249        | 1252         | rBV      | 87771          | 126422        | 6.14%           | 0.480%        |
| 19        | 9.504       | 1258          | 1261        | 1263         | rBV      | 125388         | 109369        | 5.32%           | 0.415%        |
| 20        | 9.528       | 1263          | 1265        | 1267         | rVV      | 80485          | 62493         | 3.04%           | 0.237%        |
| 21        | 9.557       | 1267          | 1270        | 1274         | rVV2     | 164074         | 184528        | 8.97%           | 0.700%        |
| 22        | 9.622       | 1278          | 1281        | 1286         | rVB      | 66894          | 71681         | 3.48%           | 0.272%        |
| 23        | 9.704       | 1292          | 1295        | 1299         | rBV      | 190973         | 163866        | 7.96%           | 0.622%        |
| 24        | 9.775       | 1305          | 1307        | 1309         | rVB      | 53675          | 34492         | 1.68%           | 0.131%        |
| 25        | 9.845       | 1312          | 1319        | 1322         | rBV      | 1327276        | 1149829       | 55.88%          | 4.362%        |
| 26        | 9.875       | 1322          | 1324        | 1328         | rVV      | 237994         | 196914        | 9.57%           | 0.747%        |
| 27        | 9.910       | 1328          | 1330        | 1333         | rVB4     | 30974          | 27567         | 1.34%           | 0.105%        |
| 28        | 9.951       | 1333          | 1337        | 1341         | rVB2     | 46863          | 60099         | 2.92%           | 0.228%        |
| 29        | 9.998       | 1341          | 1345        | 1347         | rBV4     | 28002          | 40662         | 1.98%           | 0.154%        |
| 30        | 10.045      | 1349          | 1353        | 1358         | rVV      | 164868         | 157114        | 7.64%           | 0.596%        |
| 31        | 10.086      | 1358          | 1360        | 1364         | rVV2     | 69015          | 55785         | 2.71%           | 0.212%        |
| 32        | 10.157      | 1369          | 1372        | 1375         | rBV2     | 72224          | 85379         | 4.15%           | 0.324%        |
| 33        | 10.186      | 1375          | 1377        | 1380         | rVV      | 48976          | 40113         | 1.95%           | 0.152%        |
| 34        | 10.251      | 1384          | 1388        | 1390         | rBV2     | 59185          | 78515         | 3.82%           | 0.298%        |

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Integration Parameters: rteint.p  
 Integrator: RTE  
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Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

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

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 10.275 | 1390 | 1392 | 1394 | rVB2 | 78094   | 59343   | 2.88%  | 0.225% |
| 36 | 10.298 | 1394 | 1396 | 1399 | rBV  | 33113   | 28044   | 1.36%  | 0.106% |
| 37 | 10.392 | 1408 | 1412 | 1418 | rVB  | 279953  | 279817  | 13.60% | 1.062% |
| 38 | 10.457 | 1418 | 1423 | 1425 | rBV  | 52528   | 67939   | 3.30%  | 0.258% |
| 39 | 10.498 | 1428 | 1430 | 1433 | rVV2 | 67304   | 71426   | 3.47%  | 0.271% |
| 40 | 10.545 | 1436 | 1438 | 1442 | rVB2 | 62350   | 61357   | 2.98%  | 0.233% |
| 41 | 10.633 | 1447 | 1453 | 1456 | rBV  | 368869  | 391562  | 19.03% | 1.485% |
| 42 | 10.757 | 1469 | 1474 | 1481 | rVB2 | 135391  | 187938  | 9.13%  | 0.713% |
| 43 | 10.939 | 1501 | 1505 | 1507 | rBV2 | 85504   | 99415   | 4.83%  | 0.377% |
| 44 | 10.969 | 1507 | 1510 | 1513 | rVV2 | 66788   | 73877   | 3.59%  | 0.280% |
| 45 | 11.022 | 1517 | 1519 | 1522 | rVB2 | 53817   | 48647   | 2.36%  | 0.185% |
| 46 | 11.069 | 1522 | 1527 | 1529 | rBV3 | 43532   | 66159   | 3.22%  | 0.251% |
| 47 | 11.092 | 1529 | 1531 | 1536 | rVV2 | 48825   | 55957   | 2.72%  | 0.212% |
| 48 | 11.133 | 1536 | 1538 | 1542 | rVV3 | 48654   | 50309   | 2.44%  | 0.191% |
| 49 | 11.192 | 1543 | 1548 | 1551 | rVV3 | 70582   | 113286  | 5.51%  | 0.430% |
| 50 | 11.227 | 1551 | 1554 | 1560 | rVB  | 168218  | 169423  | 8.23%  | 0.643% |
| 51 | 11.327 | 1568 | 1571 | 1573 | rBV  | 1123712 | 1021868 | 49.66% | 3.877% |
| 52 | 11.351 | 1573 | 1575 | 1579 | rVV  | 1431974 | 1268515 | 61.65% | 4.812% |
| 53 | 11.404 | 1581 | 1584 | 1587 | rVB  | 723015  | 571652  | 27.78% | 2.169% |
| 54 | 11.439 | 1587 | 1590 | 1593 | rBV2 | 69703   | 75793   | 3.68%  | 0.288% |
| 55 | 11.563 | 1608 | 1611 | 1616 | rBV  | 159278  | 156980  | 7.63%  | 0.596% |
| 56 | 11.622 | 1618 | 1621 | 1624 | rVB  | 96382   | 76153   | 3.70%  | 0.289% |
| 57 | 11.663 | 1625 | 1628 | 1633 | rVB2 | 68863   | 79324   | 3.85%  | 0.301% |
| 58 | 11.722 | 1636 | 1638 | 1640 | rBV  | 39417   | 32792   | 1.59%  | 0.124% |
| 59 | 11.739 | 1640 | 1641 | 1646 | rVB  | 48912   | 47384   | 2.30%  | 0.180% |
| 60 | 11.786 | 1646 | 1649 | 1651 | rBV2 | 30742   | 32564   | 1.58%  | 0.124% |
| 61 | 11.833 | 1653 | 1657 | 1659 | rBV  | 217572  | 203438  | 9.89%  | 0.772% |
| 62 | 11.857 | 1659 | 1661 | 1665 | rVB  | 300650  | 258751  | 12.57% | 0.982% |
| 63 | 11.904 | 1666 | 1669 | 1672 | rBV  | 142212  | 139852  | 6.80%  | 0.531% |
| 64 | 11.945 | 1672 | 1676 | 1678 | rVV2 | 414048  | 448260  | 21.78% | 1.700% |
| 65 | 11.963 | 1678 | 1679 | 1684 | rVB  | 178270  | 143141  | 6.96%  | 0.543% |
| 66 | 12.033 | 1688 | 1691 | 1693 | rVV2 | 61768   | 62490   | 3.04%  | 0.237% |
| 67 | 12.063 | 1694 | 1696 | 1699 | rVB2 | 42491   | 34555   | 1.68%  | 0.131% |
| 68 | 12.127 | 1704 | 1707 | 1709 | rVV  | 167825  | 149727  | 7.28%  | 0.568% |
| 69 | 12.151 | 1709 | 1711 | 1715 | rVB2 | 60712   | 72306   | 3.51%  | 0.274% |
| 70 | 12.257 | 1727 | 1729 | 1730 | rBV2 | 42072   | 34178   | 1.66%  | 0.130% |
| 71 | 12.274 | 1730 | 1732 | 1735 | rVB  | 68036   | 49482   | 2.40%  | 0.188% |

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Integration Parameters: rteint.p  
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Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF071620.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 12.304 | 1735 | 1737 | 1740 | rBV  | 105664  | 101538  | 4.93%   | 0.385% |
| 73  | 12.380 | 1747 | 1750 | 1753 | rBV  | 191806  | 200313  | 9.73%   | 0.760% |
| 74  | 12.421 | 1753 | 1757 | 1762 | rVV5 | 207957  | 392256  | 19.06%  | 1.488% |
| 75  | 12.469 | 1762 | 1765 | 1769 | rVV2 | 112784  | 126396  | 6.14%   | 0.479% |
| 76  | 12.545 | 1774 | 1778 | 1781 | rVV  | 2137794 | 1778905 | 86.45%  | 6.748% |
| 77  | 12.586 | 1783 | 1785 | 1787 | rVV  | 57316   | 54022   | 2.63%   | 0.205% |
| 78  | 12.639 | 1791 | 1794 | 1796 | rBV2 | 87141   | 79719   | 3.87%   | 0.302% |
| 79  | 12.698 | 1801 | 1804 | 1807 | rVV3 | 83923   | 107736  | 5.24%   | 0.409% |
| 80  | 12.774 | 1811 | 1817 | 1820 | rBV  | 1908815 | 2057695 | 100.00% | 7.806% |
| 81  | 12.827 | 1823 | 1826 | 1829 | rVB2 | 77223   | 104821  | 5.09%   | 0.398% |
| 82  | 12.892 | 1834 | 1837 | 1838 | rBV2 | 102534  | 99054   | 4.81%   | 0.376% |
| 83  | 12.910 | 1838 | 1840 | 1848 | rVB  | 708695  | 721991  | 35.09%  | 2.739% |
| 84  | 12.986 | 1850 | 1853 | 1857 | rBV2 | 121463  | 178303  | 8.67%   | 0.676% |
| 85  | 13.074 | 1866 | 1868 | 1869 | rBV2 | 90195   | 73459   | 3.57%   | 0.279% |
| 86  | 13.098 | 1869 | 1872 | 1875 | rVB  | 350231  | 353043  | 17.16%  | 1.339% |
| 87  | 13.145 | 1878 | 1880 | 1881 | rBV  | 71073   | 54984   | 2.67%   | 0.209% |
| 88  | 13.163 | 1881 | 1883 | 1886 | rVB  | 268289  | 241161  | 11.72%  | 0.915% |
| 89  | 13.204 | 1887 | 1890 | 1893 | rBV2 | 206924  | 190101  | 9.24%   | 0.721% |
| 90  | 13.298 | 1903 | 1906 | 1908 | rVB  | 115413  | 88006   | 4.28%   | 0.334% |
| 91  | 13.327 | 1908 | 1911 | 1914 | rBV  | 136421  | 125796  | 6.11%   | 0.477% |
| 92  | 13.745 | 1980 | 1982 | 1984 | rBV  | 109153  | 74348   | 3.61%   | 0.282% |
| 93  | 13.768 | 1984 | 1986 | 1988 | rBV  | 96917   | 97004   | 4.71%   | 0.368% |
| 94  | 13.816 | 1992 | 1994 | 1998 | rVB2 | 154014  | 152827  | 7.43%   | 0.580% |
| 95  | 13.968 | 2015 | 2020 | 2023 | rBV2 | 1402695 | 1943219 | 94.44%  | 7.372% |
| 96  | 13.998 | 2023 | 2025 | 2028 | rVB  | 779148  | 585288  | 28.44%  | 2.220% |
| 97  | 14.074 | 2036 | 2038 | 2041 | rBV  | 174050  | 134356  | 6.53%   | 0.510% |
| 98  | 15.004 | 2193 | 2196 | 2203 | rVB  | 518962  | 966378  | 46.96%  | 3.666% |
| 99  | 15.362 | 2254 | 2257 | 2263 | rVB  | 451983  | 534850  | 25.99%  | 2.029% |
| 100 | 15.427 | 2264 | 2268 | 2271 | rBV  | 567300  | 773314  | 37.58%  | 2.934% |

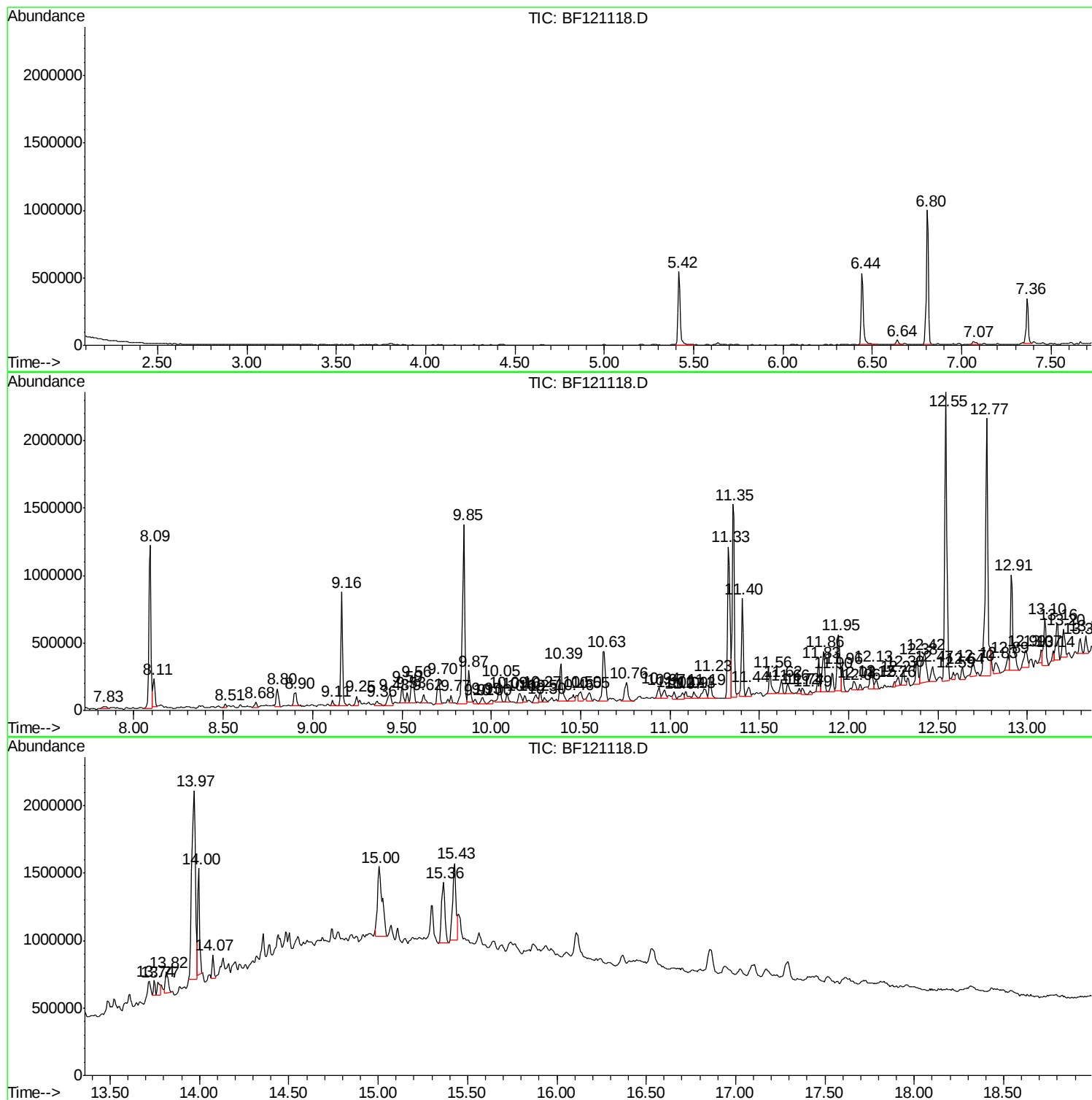
Sum of corrected areas: 26360497

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

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
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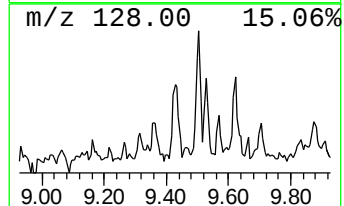
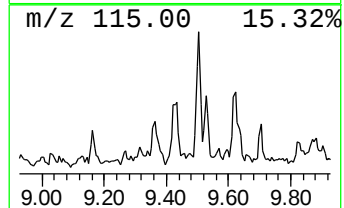
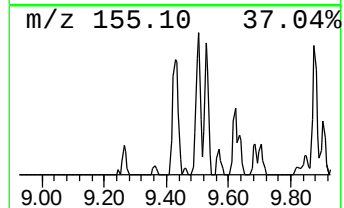
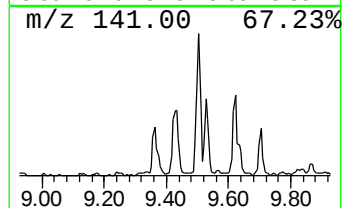
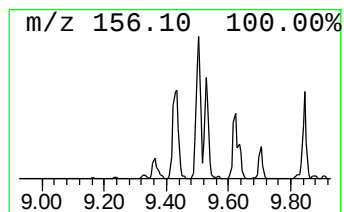
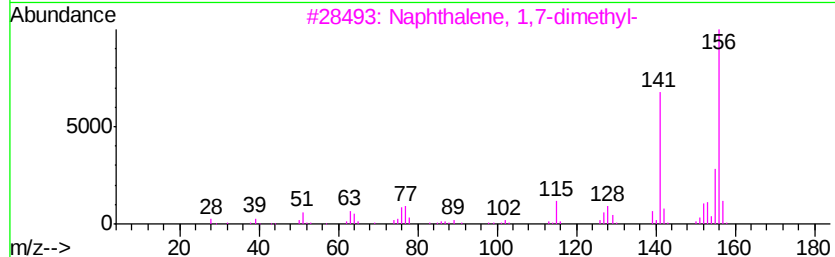
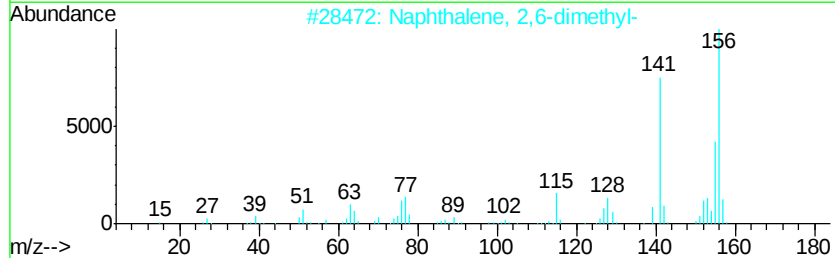
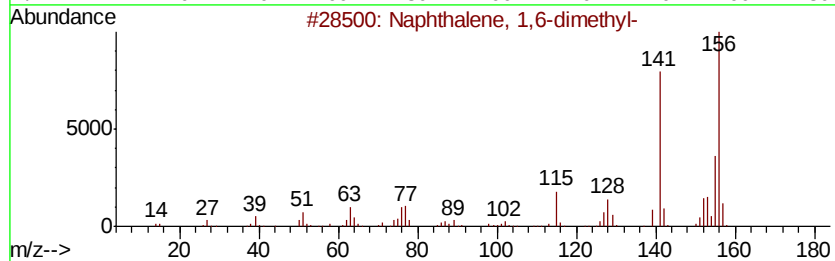
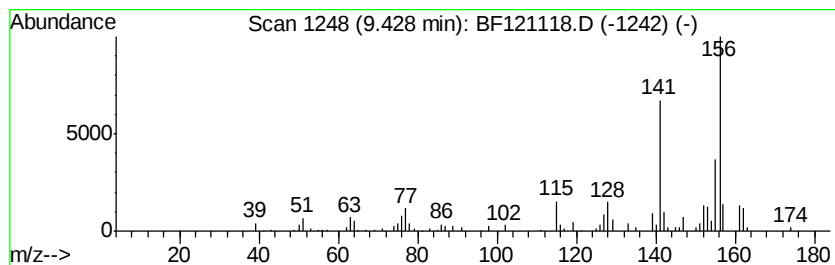
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Naphthalene, 1,6-dimethyl- Concentration Rank 16

| R.T.    | EstConc | Area                       | Relative to ISTD | R.T.           |
|---------|---------|----------------------------|------------------|----------------|
| 9.43    | 2.20 ng | 126422                     | Acenaphthene-d10 | 9.85           |
| Hit# of | 5       | Tentative ID               | MW MolForm       | CAS# Qual      |
| 1       |         | Naphthalene, 1,6-dimethyl- | 156 C12H12       | 000575-43-9 97 |
| 2       |         | Naphthalene, 2,6-dimethyl- | 156 C12H12       | 000581-42-0 97 |
| 3       |         | Naphthalene, 1,7-dimethyl- | 156 C12H12       | 000575-37-1 97 |
| 4       |         | Naphthalene, 2,3-dimethyl- | 156 C12H12       | 000581-40-8 96 |
| 5       |         | Naphthalene, 2,7-dimethyl- | 156 C12H12       | 000582-16-1 96 |



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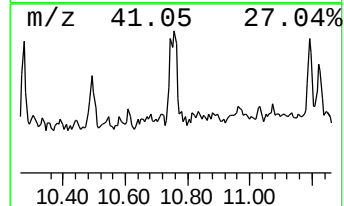
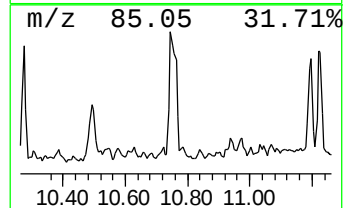
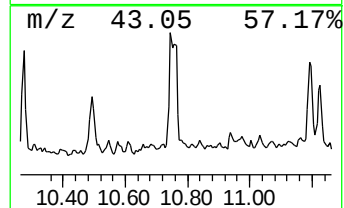
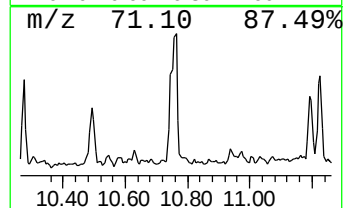
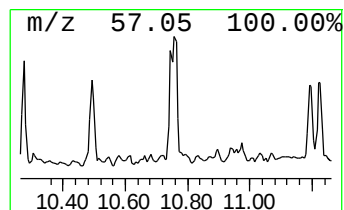
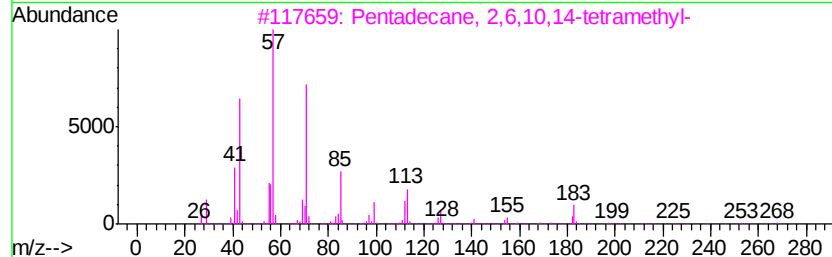
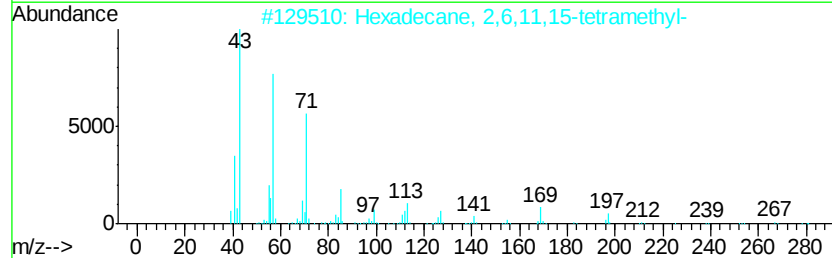
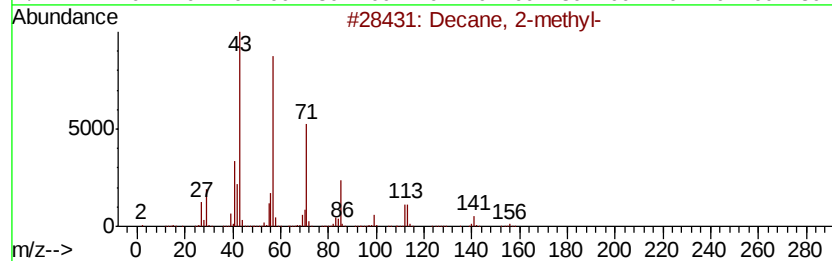
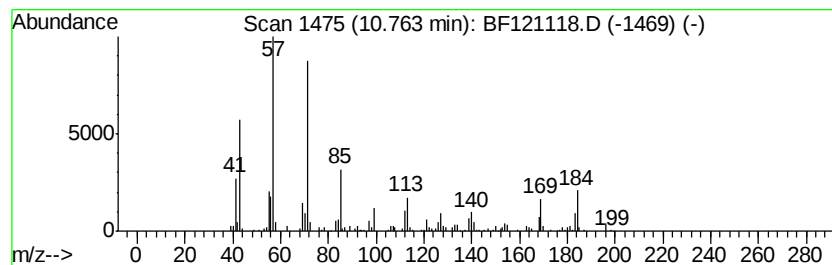
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 unknown10.76 Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 10.76     | 3.68 ng                             | 187938 | Phenanthrene-d10 | 11.33       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Decane, 2-methyl-                   | 156    | C11H24           | 006975-98-0 | 46   |
| 2         | Hexadecane, 2,6,11,15-tetramethyl-  | 282    | C20H42           | 000504-44-9 | 46   |
| 3         | Pentadecane, 2,6,10,14-tetramethyl- | 268    | C19H40           | 001921-70-6 | 45   |
| 4         | Hexadecane, 7,9-dimethyl-           | 254    | C18H38           | 021164-95-4 | 43   |
| 5         | Octadecane, 2,6-dimethyl-           | 282    | C20H42           | 075163-97-2 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

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

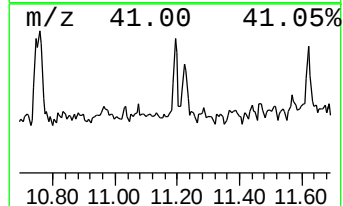
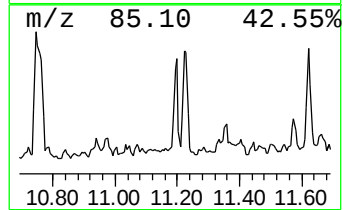
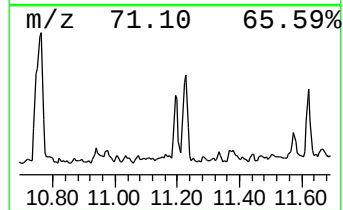
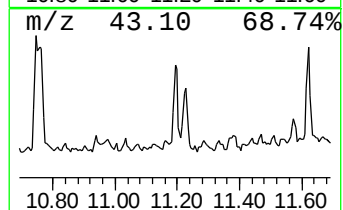
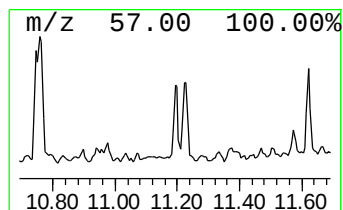
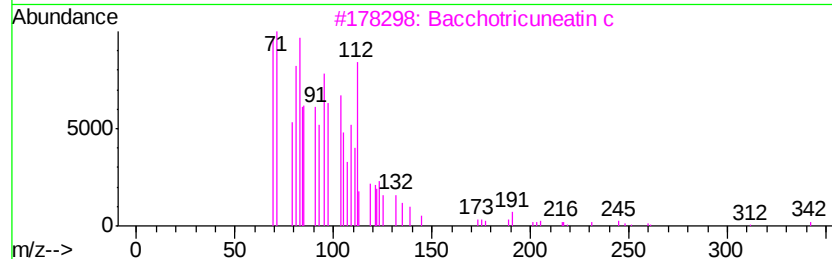
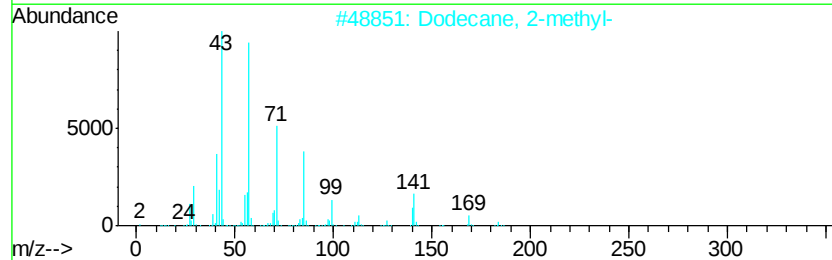
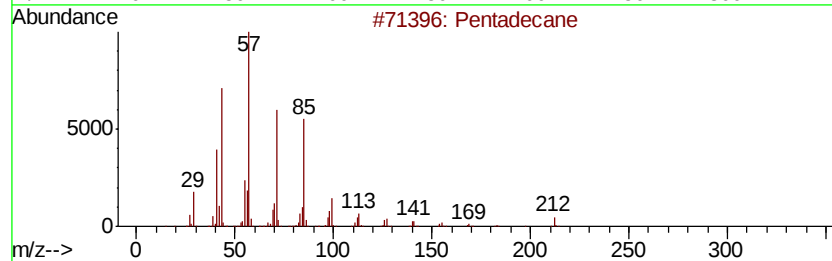
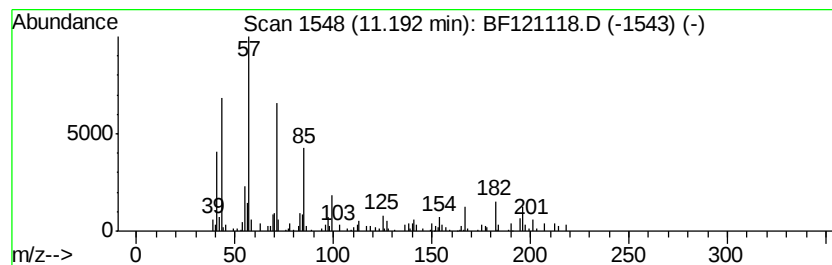
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 Pentadecane Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.19 | 2.22 ng | 113286 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | Pentadecane         | 212 | C15H32   | 000629-62-9 | 89   |
| 2    |    | Dodecane, 2-methyl- | 184 | C13H28   | 001560-97-0 | 62   |
| 3    |    | Bacchotricuneatin c | 342 | C20H22O5 | 066563-30-2 | 60   |
| 4    |    | Tetracosane         | 338 | C24H50   | 000646-31-1 | 58   |
| 5    |    | Heptacosane         | 380 | C27H56   | 000593-49-7 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

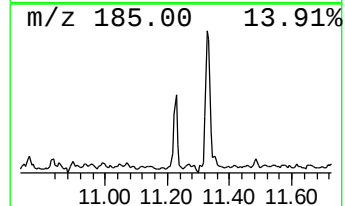
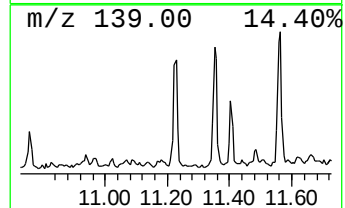
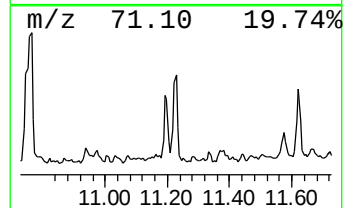
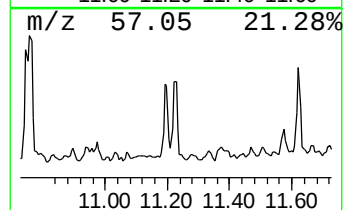
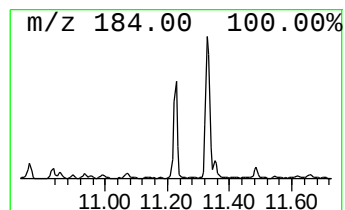
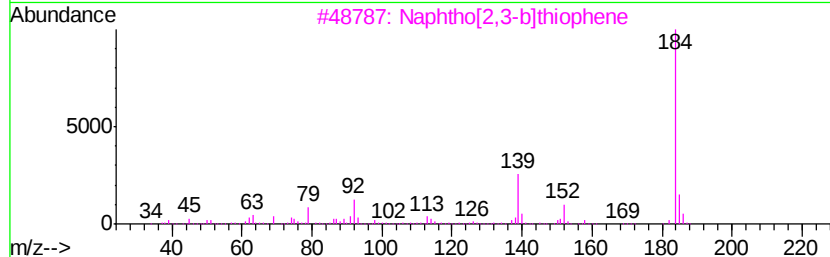
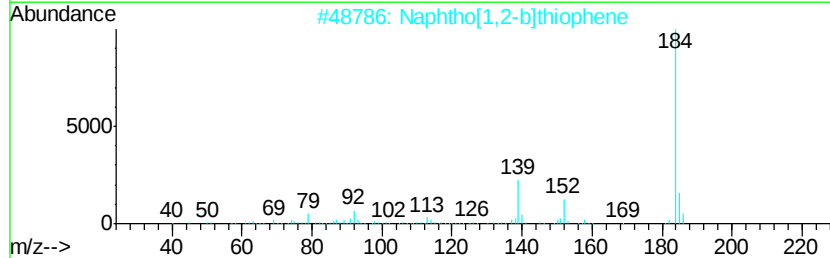
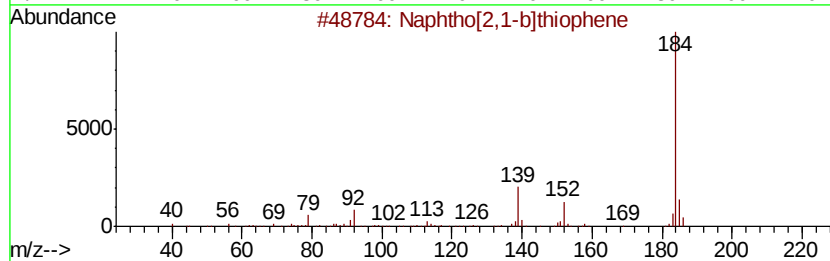
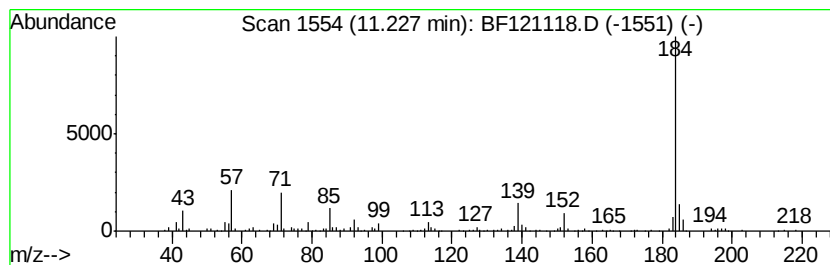
Instrument :  
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ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Naphtho[2,1-b]thiophene Concentration Rank 8

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.    |             |      |
|---------|---------|-------------------------|------------------|---------|-------------|------|
| 11.23   | 3.32 ng | 169423                  | Phenanthrene-d10 | 11.33   |             |      |
| Hit# of | 5       | Tentative ID            | MW               | MolForm | CAS#        | Qual |
| 1       |         | Naphtho[2,1-b]thiophene | 184              | C12H8S  | 000233-02-3 | 93   |
| 2       |         | Naphtho[1,2-b]thiophene | 184              | C12H8S  | 000234-41-3 | 93   |
| 3       |         | Naphtho[2,3-b]thiophene | 184              | C12H8S  | 000268-77-9 | 93   |
| 4       |         | Dibenzothiophene        | 184              | C12H8S  | 000132-65-0 | 89   |
| 5       |         | Azuleno(2,1-b)thiophene | 184              | C12H8S  | 000248-13-5 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

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

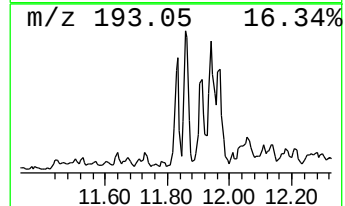
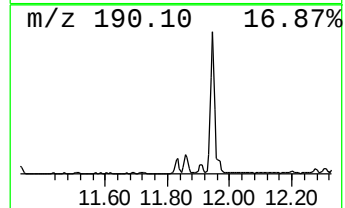
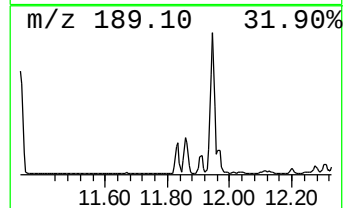
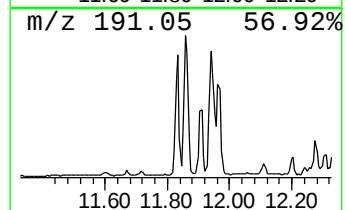
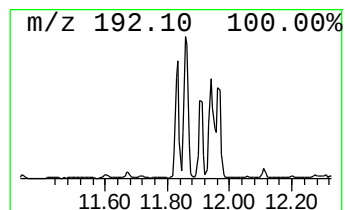
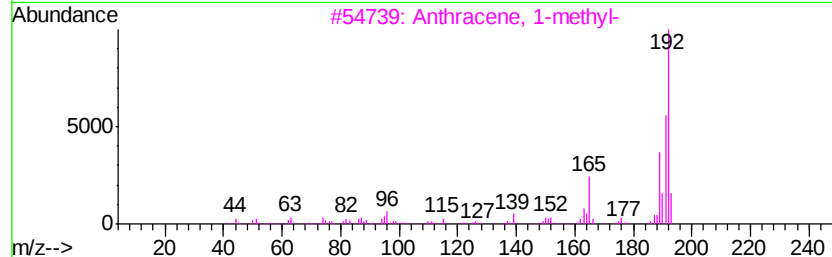
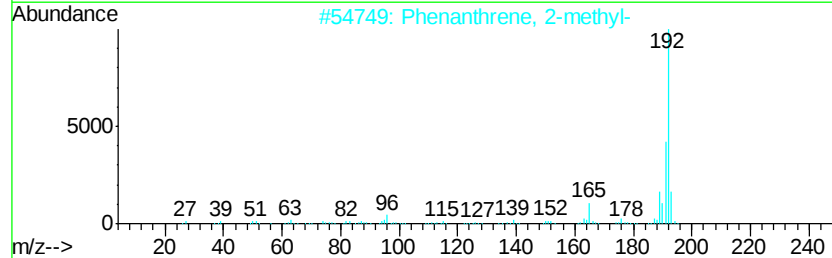
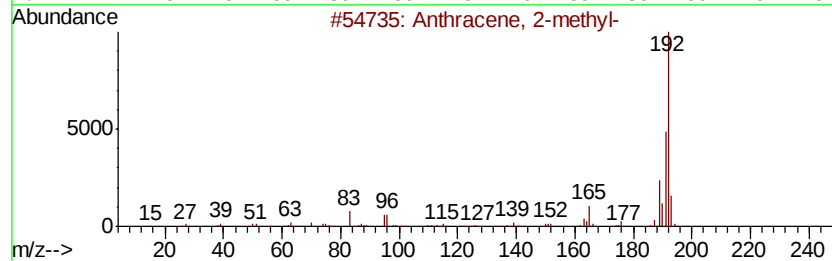
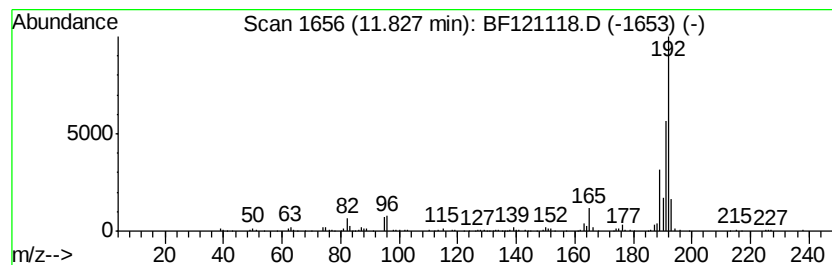
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Anthracene, 2-methyl- Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.83 | 3.98 ng | 203438 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 96   |
| 2    |    | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 95   |
| 3    |    | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 95   |
| 4    |    | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 94   |
| 5    |    | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

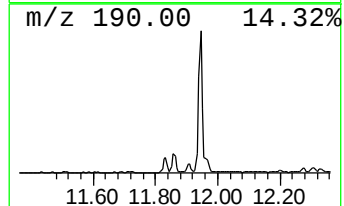
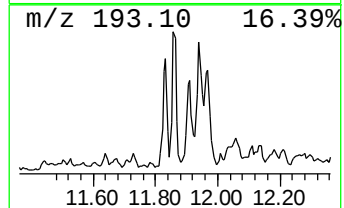
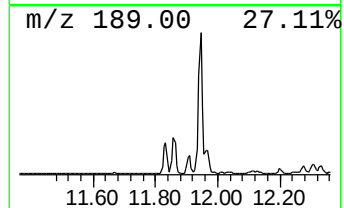
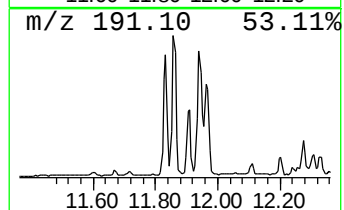
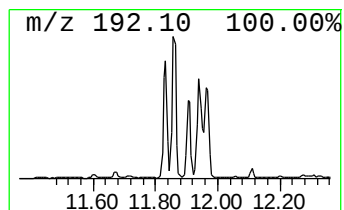
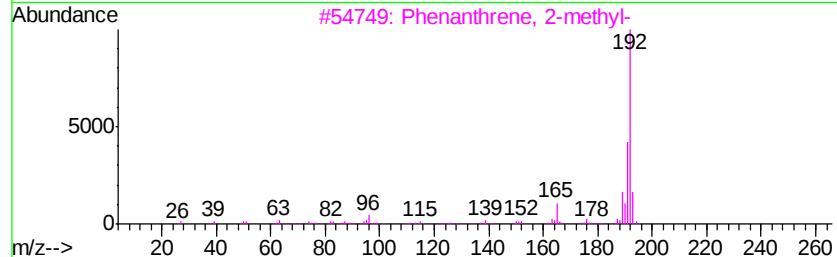
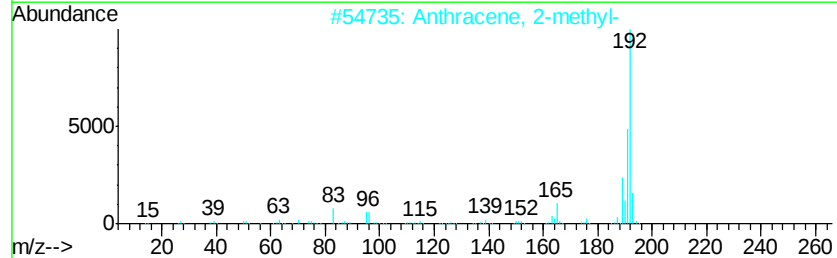
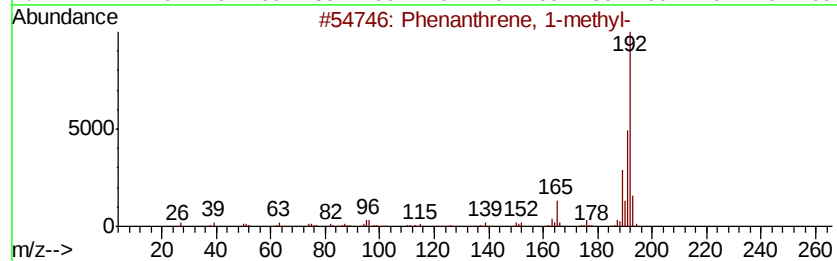
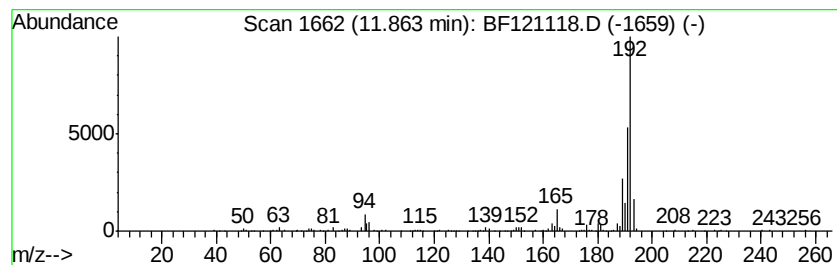
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ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Phenanthrene, 1-methyl- Concentration Rank 3

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.86     | 5.06 ng                             | 258751 | Phenanthrene-d10 | 11.33       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 97   |
| 2         | Anthracene, 2-methyl-               | 192    | C15H12           | 000613-12-7 | 97   |
| 3         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 97   |
| 4         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 96   |
| 5         | Anthracene, 9-methyl-               | 192    | C15H12           | 000779-02-2 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
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Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

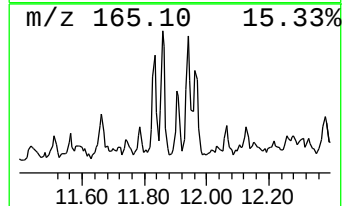
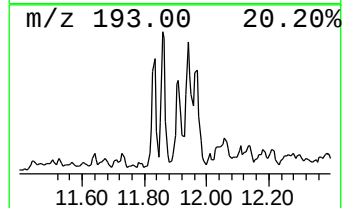
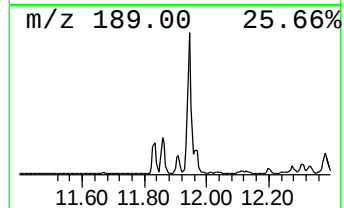
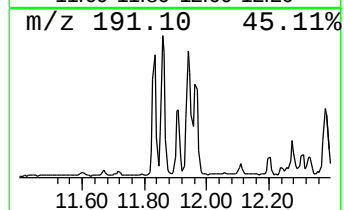
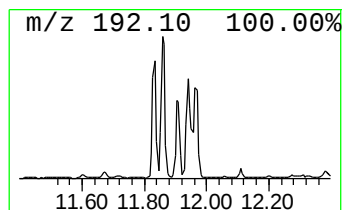
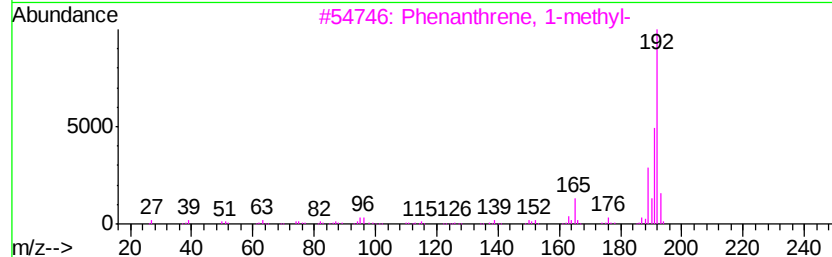
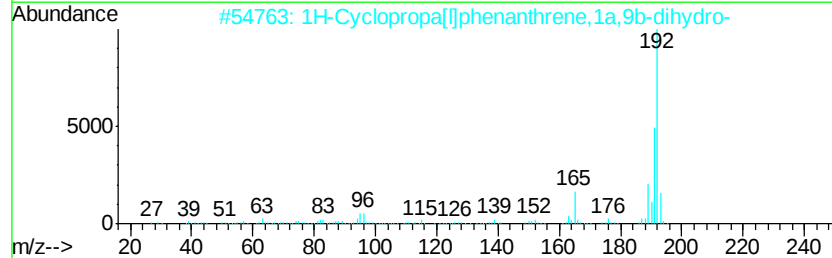
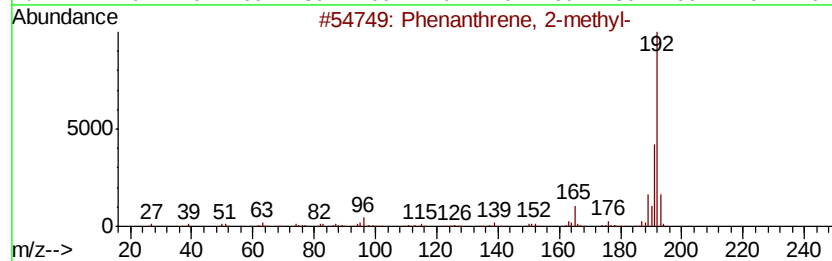
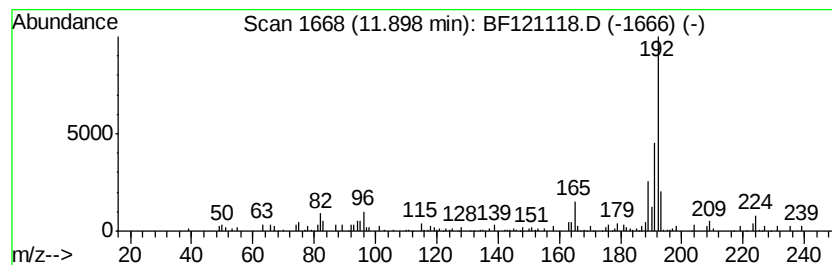
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 11

| R.T.      | EstConc                             | Area   | Relative to ISTD |             | R.T.  |
|-----------|-------------------------------------|--------|------------------|-------------|-------|
| 11.90     | 2.74 ng                             | 139852 | Phenanthrene-d10 |             | 11.33 |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual  |
| 1         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 97    |
| 2         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 97    |
| 3         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 96    |
| 4         | Anthracene, 2-methyl-               | 192    | C15H12           | 000613-12-7 | 96    |
| 5         | 1H-Indene, 2-phenyl-                | 192    | C15H12           | 004505-48-0 | 95    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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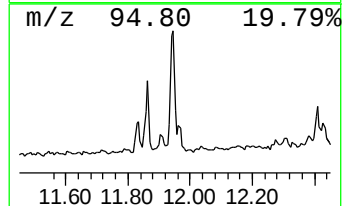
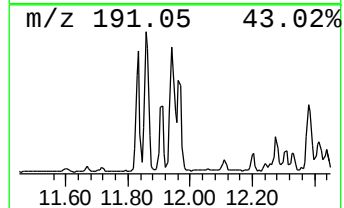
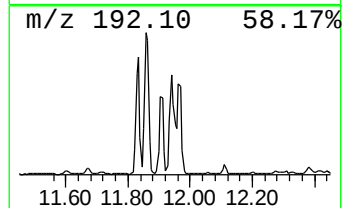
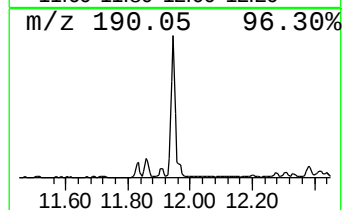
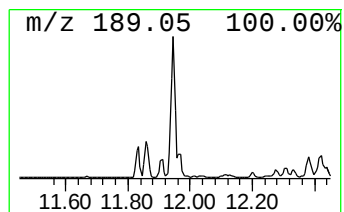
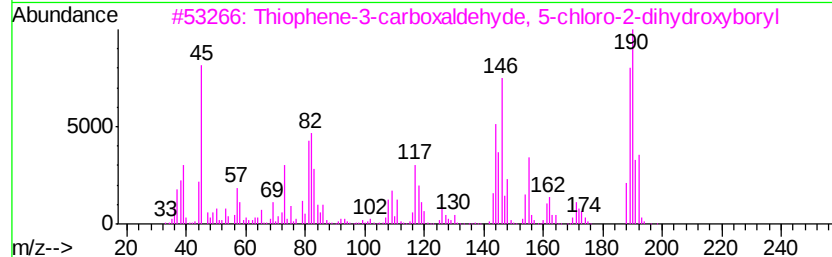
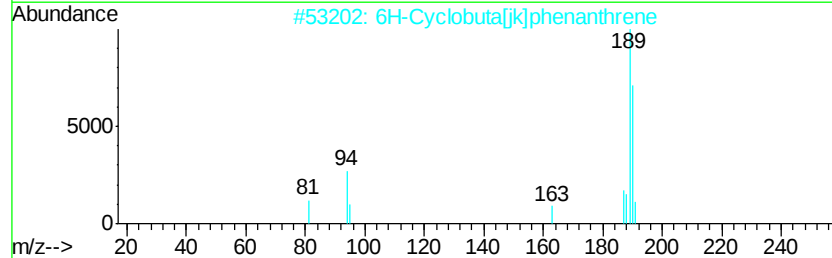
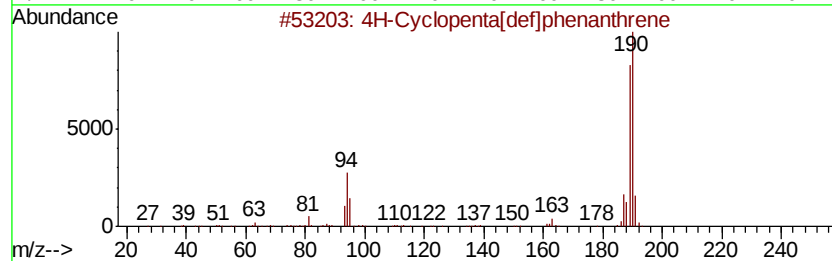
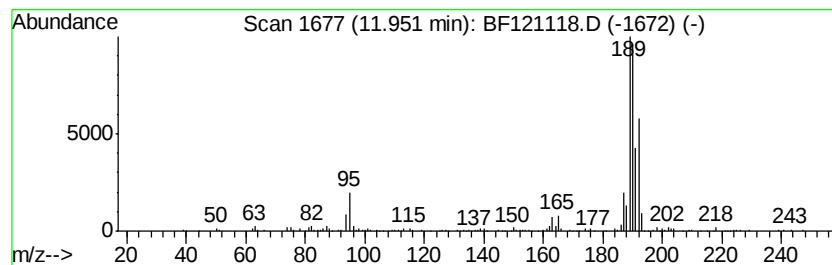
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.95 | 8.77 ng | 448260 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                                          | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------------------------|-----|------------|-------------|------|
| 1    | 5  | 4H-Cyclopenta[def]phenanthrene                        | 190 | C15H10     | 000203-64-5 | 70   |
| 2    |    | 6H-Cyclobuta[ik]phenanthrene                          | 190 | C15H10     | 083469-43-6 | 59   |
| 3    |    | Thiophene-3-carboxaldehyde, 5-chloro-2-dihydroxyboryl | 190 | C5H4BClO3S | 036155-87-0 | 52   |
| 4    |    | 2,2'-Bis(4,5-dimethylimidazole)                       | 190 | C10H14N4   | 069286-06-2 | 42   |
| 5    |    | 1,4-Naphthalenedione, 2,5-dihydroxy                   | 190 | C10H6O4    | 004923-55-1 | 27   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

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

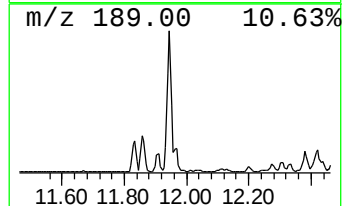
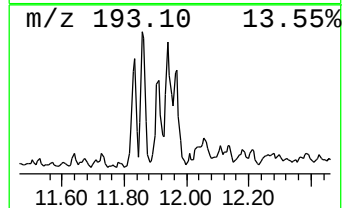
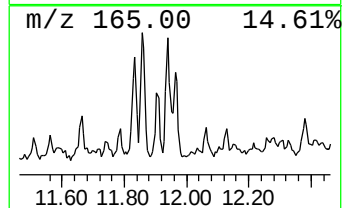
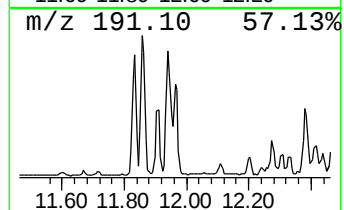
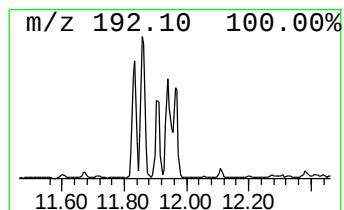
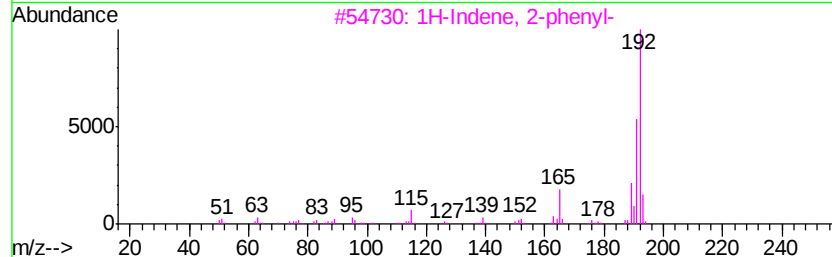
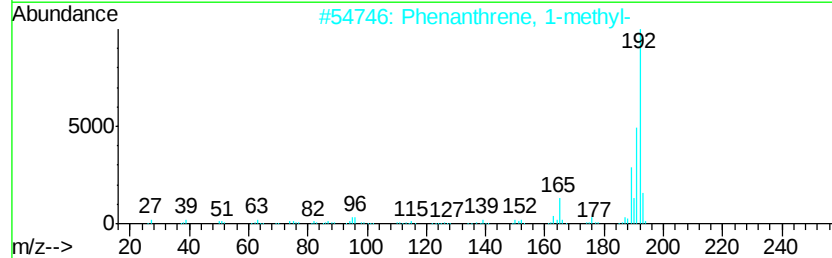
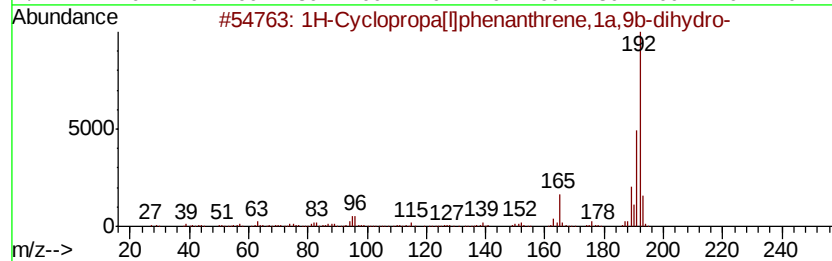
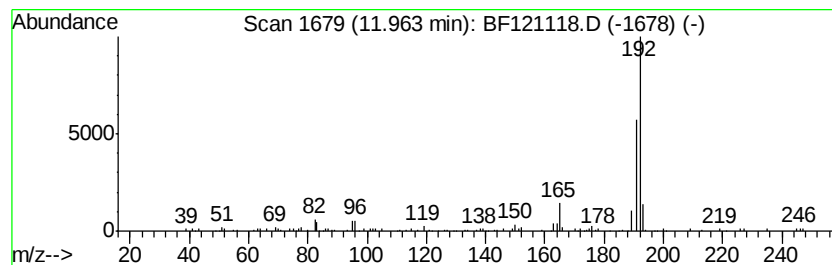
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 1H-Cyclopropa[l]phenanthren... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.96 | 2.80 ng | 143141 | Phenanthrene-d10 | 11.33 |

| Hit# | of | 5 | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Cyclopropa[l]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 91   |
| 2    |    |   | Phenanthrene, 1-methyl-               | 192 | C15H12  | 000832-69-9 | 87   |
| 3    |    |   | 1H-Indene, 2-phenyl-                  | 192 | C15H12  | 004505-48-0 | 87   |
| 4    |    |   | Anthracene, 1-methyl-                 | 192 | C15H12  | 000610-48-0 | 87   |
| 5    |    |   | Anthracene, 2-methyl-                 | 192 | C15H12  | 000613-12-7 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
17

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

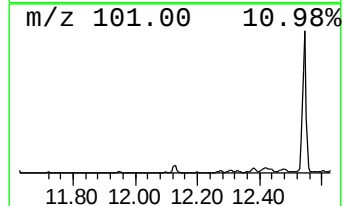
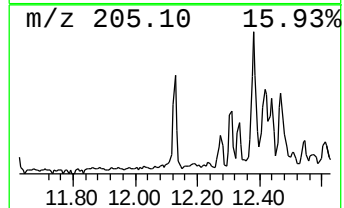
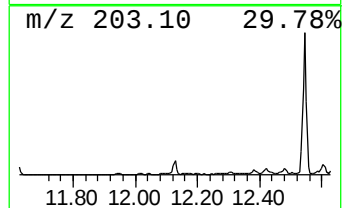
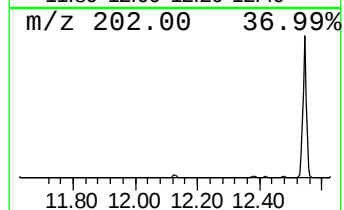
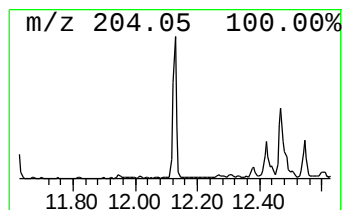
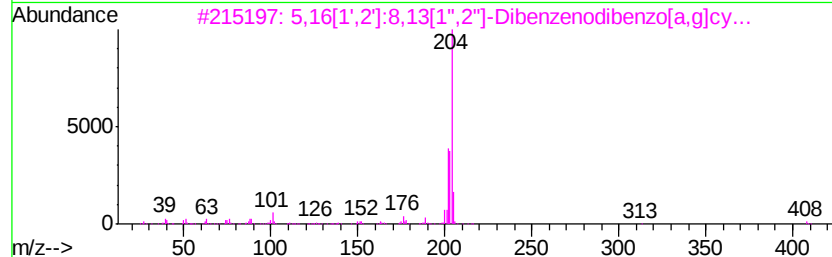
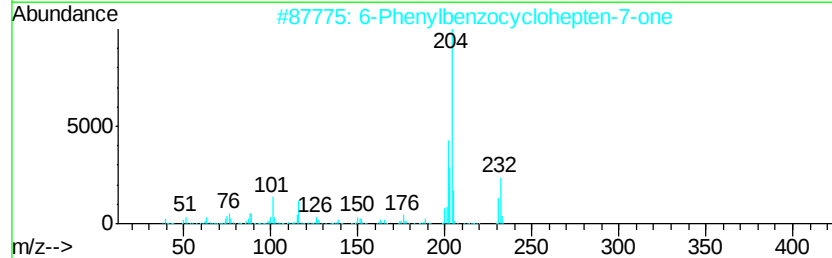
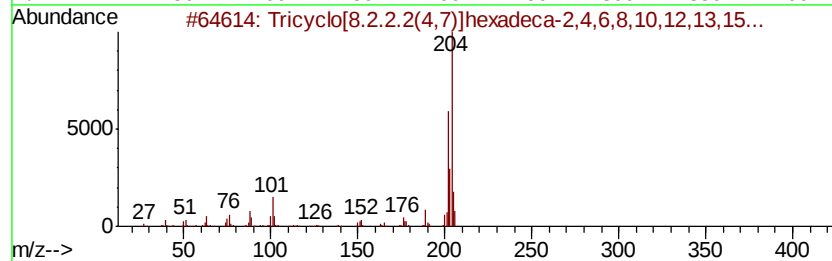
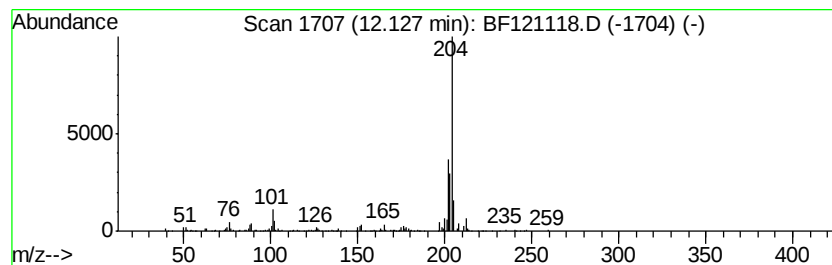
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.13 | 2.93 ng | 149727 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12  | 006572-60-7 | 90   |
| 2    |    | 6-Phenylbenzocyclohepten-7-one      | 232 | C17H12O | 093327-56-1 | 86   |
| 3    |    | 5,16[1',2':8,13[1'',2'']-Dibenz...  | 408 | C32H24  | 005672-97-9 | 86   |
| 4    |    | Naphthalene, 2-phenyl-              | 204 | C16H12  | 000612-94-2 | 76   |
| 5    |    | Naphthalene, 1-phenyl-              | 204 | C16H12  | 000605-02-7 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

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

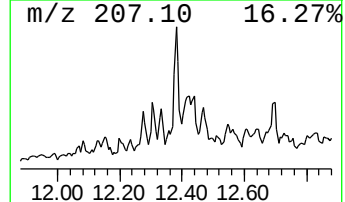
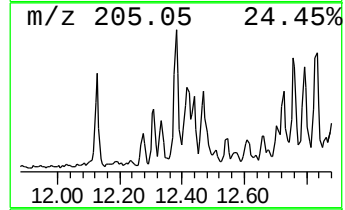
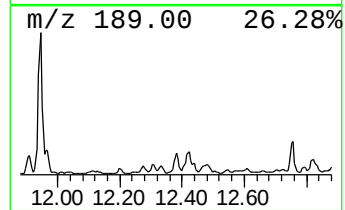
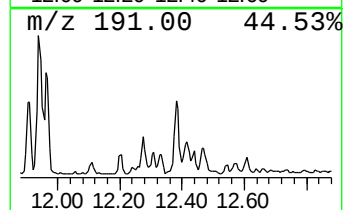
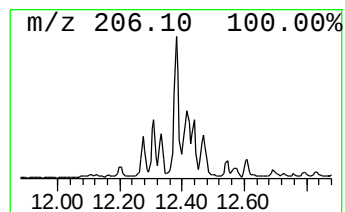
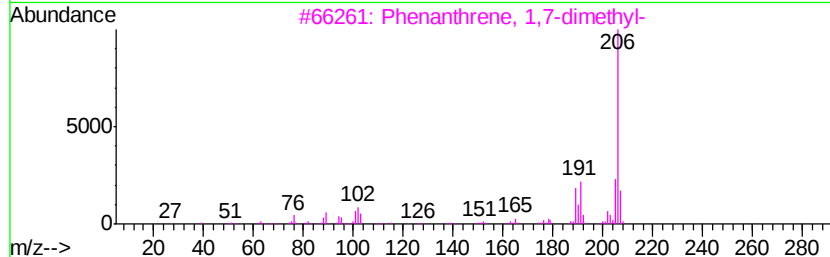
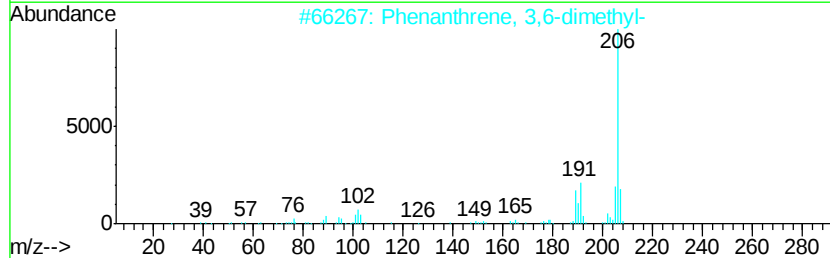
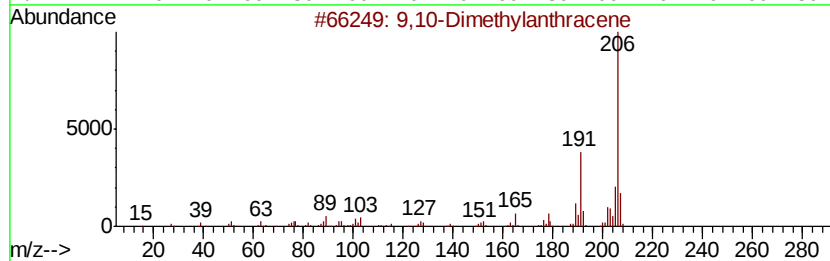
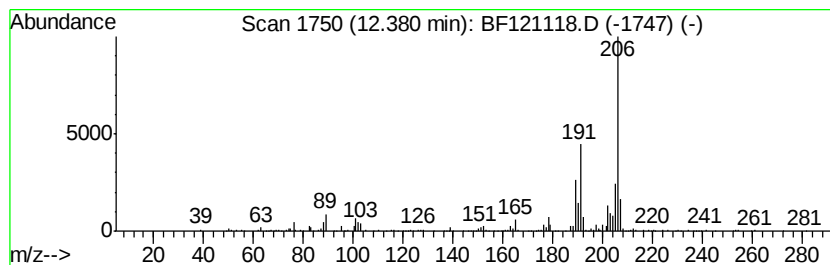
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 9,10-Dimethylantracene Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.38 | 3.92 ng | 200313 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 94   |
| 2    |    | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 94   |
| 3    |    | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 93   |
| 4    |    | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 91   |
| 5    |    | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
17

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

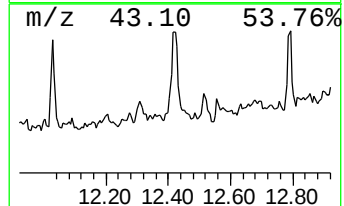
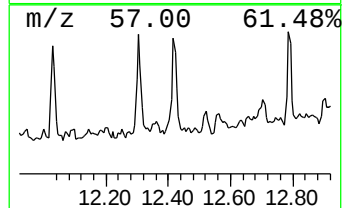
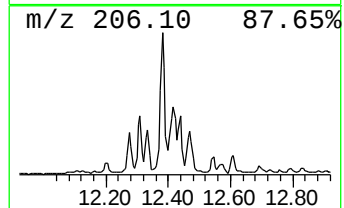
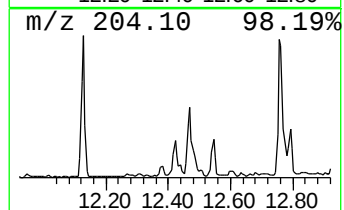
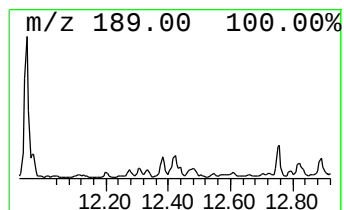
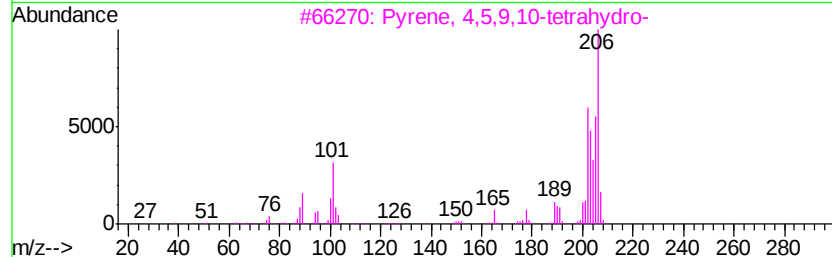
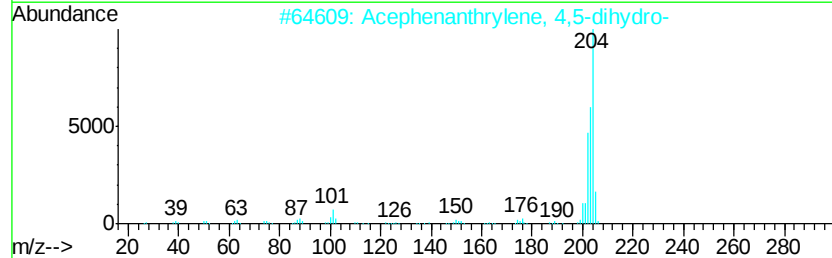
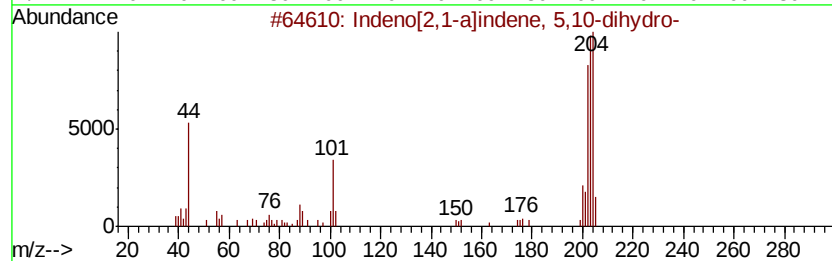
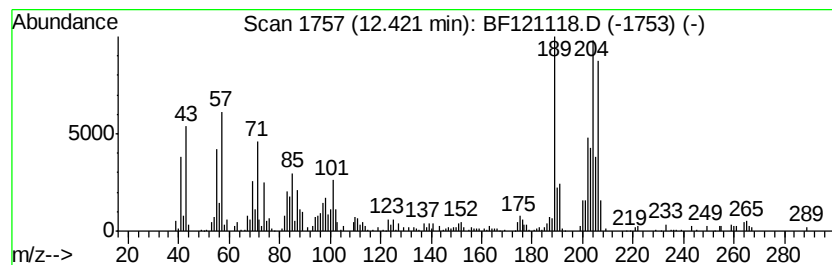
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 unknown12.42 Concentration Rank 2

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.42 | 7.68 ng | 392256 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                       | MW  | MolForm   | CAS#        | Qual |
|------|----|------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Indeno[2,1-a]indene, 5,10-dihydro- | 204 | C16H12    | 006543-29-9 | 45   |
| 2    |    | Acephenanthrylene, 4,5-dihydro-    | 204 | C16H12    | 006232-48-0 | 38   |
| 3    |    | Pvrene, 4,5,9,10-tetrahydro-       | 206 | C16H14    | 000781-17-9 | 38   |
| 4    |    | 9,10-Bis(bromomethyl)anthracene    | 362 | C16H12Br2 | 034373-96-1 | 35   |
| 5    |    | Phenanthrene, 3,6-dimethyl-        | 206 | C16H14    | 001576-67-6 | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
17

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

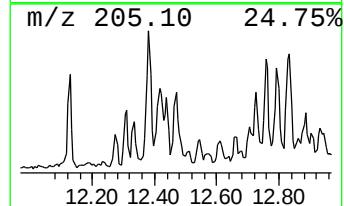
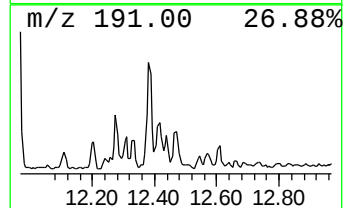
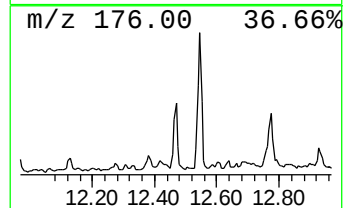
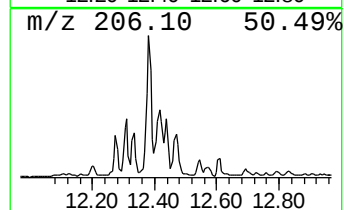
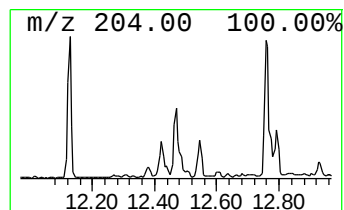
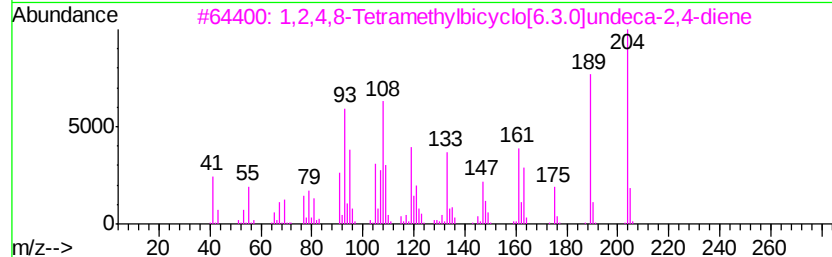
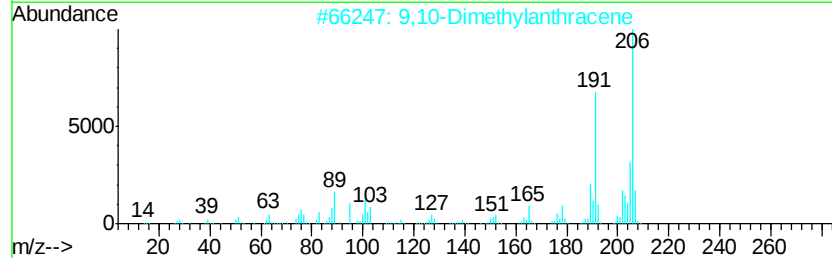
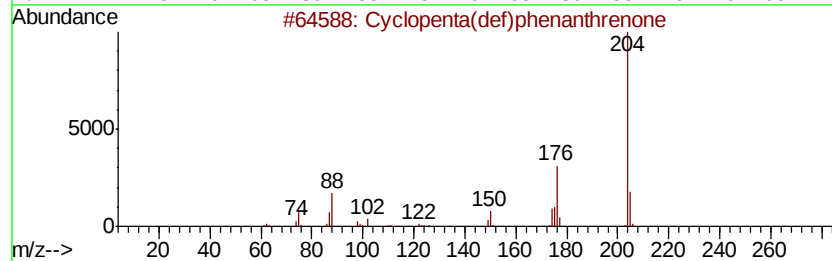
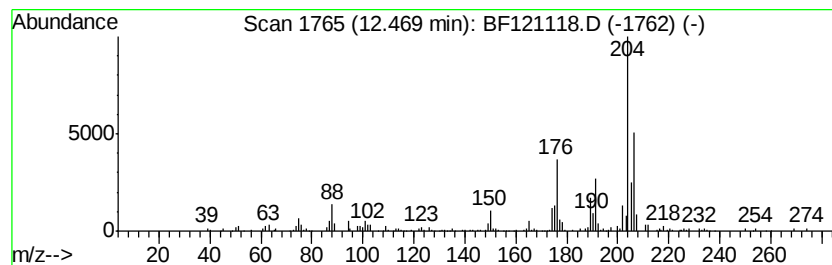
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Cyclopenta(def)phenanthrenone Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.47 | 2.47 ng | 126396 | Phenanthrene-d10 | 11.33 |

| Hit# | of | Tentative ID                                      | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrenone                     | 204 | C15H8O  | 005737-13-3 | 87   |
| 2    |    | 9,10-Dimethylantracene                            | 206 | C16H14  | 000781-43-1 | 42   |
| 3    |    | 1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene | 204 | C15H24  | 137235-51-9 | 41   |
| 4    |    | Benzene, (cyclopropylidenephenvl...               | 206 | C16H14  | 007632-57-7 | 38   |
| 5    |    | Phenanthrene, 3,6-dimethyl-                       | 206 | C16H14  | 001576-67-6 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
17

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

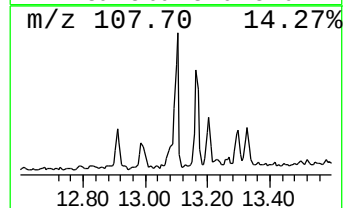
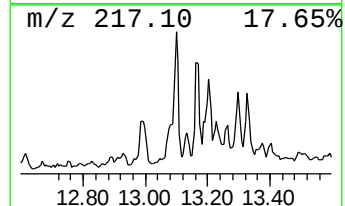
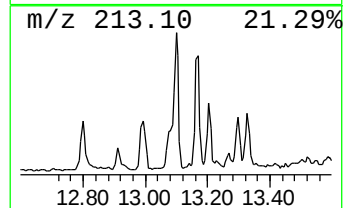
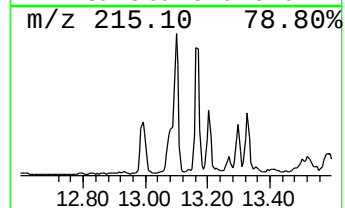
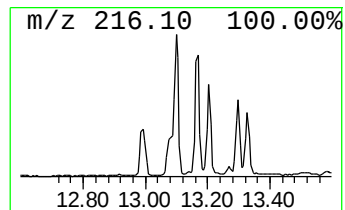
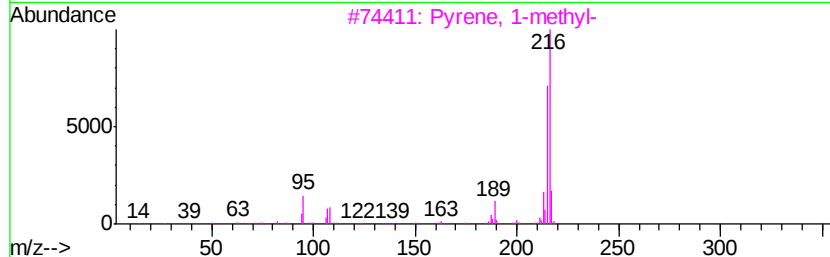
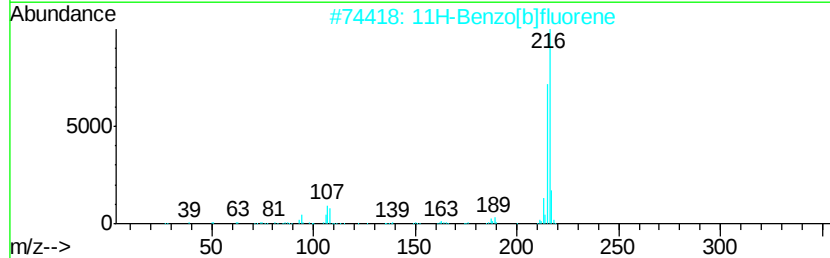
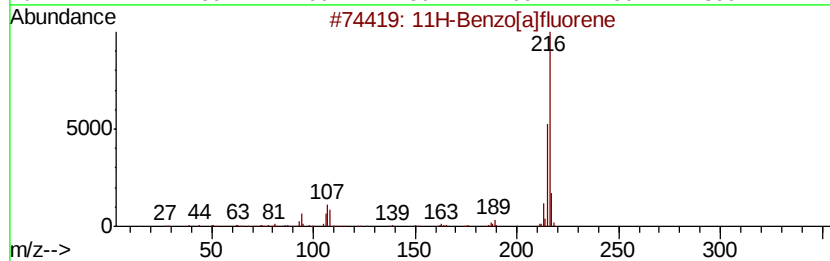
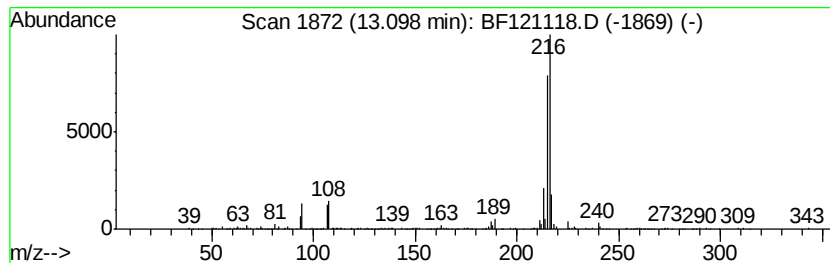
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 11H-Benzo[a]fluorene Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.10 | 3.63 ng | 353043 | Chrysene-d12     | 13.97 |

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072920\  
Data File : BF121118.D  
Acq On : 29 Jul 2020 23:33  
Operator : JU/CG  
Sample : L3436-17 5X  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
17

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

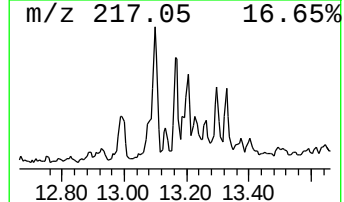
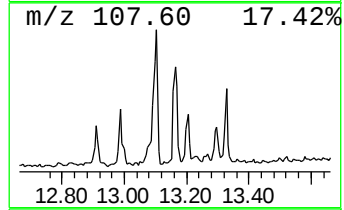
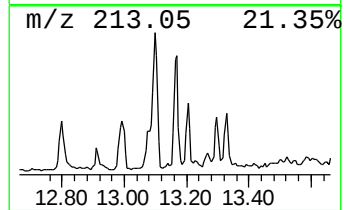
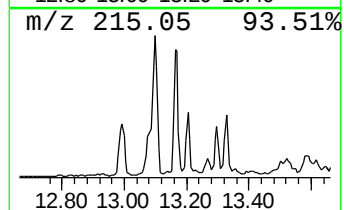
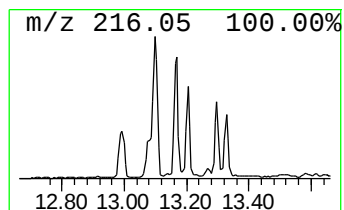
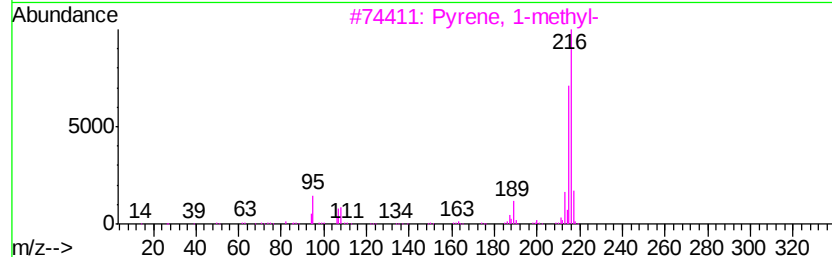
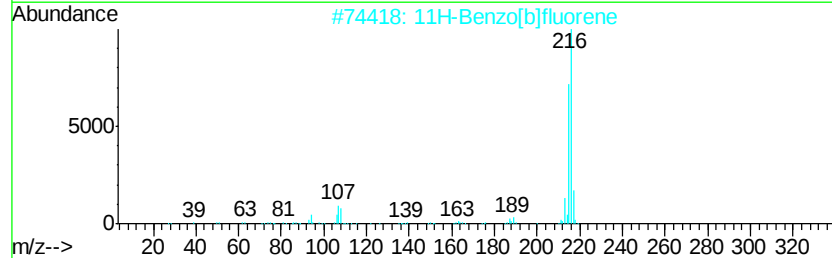
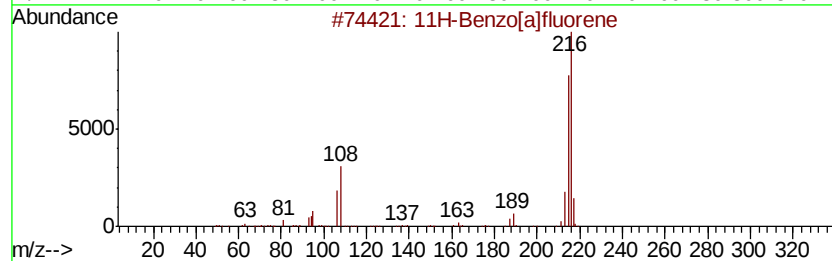
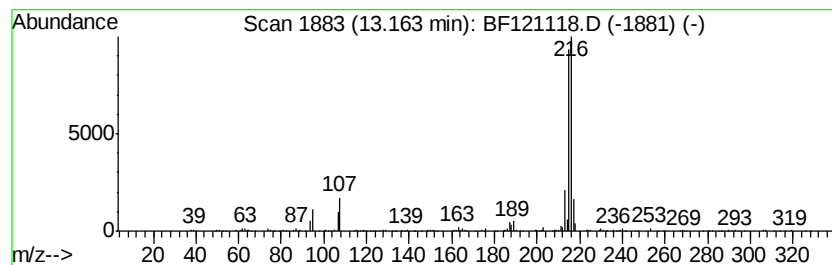
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 11H-Benzo[b]fluorene Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.16 | 2.48 ng | 241161 | Chrysene-d12     | 13.97 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 91   |
| 2    |    | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 90   |
| 3    |    | Pvrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 87   |
| 4    |    | Pvrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 76   |
| 5    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF072920\  
 Data File : BF121118.D  
 Acq On : 29 Jul 2020 23:33  
 Operator : JU/CG  
 Sample : L3436-17 5X  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 17

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

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

| TIC Top Hit name      | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|-----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                       |       |         |       |          | #                     | RT    | Resp    | Conc |
| Naphthalene, 1,6-...  | 9.43  | 2.2     | ng    | 126422   | 3                     | 9.85  | 1149830 | 20.0 |
| unknown10.76          | 10.76 | 3.7     | ng    | 187938   | 4                     | 11.33 | 1021870 | 20.0 |
| Pentadecane           | 11.19 | 2.2     | ng    | 113286   | 4                     | 11.33 | 1021870 | 20.0 |
| Naphtho[2,1-b]thi...  | 11.23 | 3.3     | ng    | 169423   | 4                     | 11.33 | 1021870 | 20.0 |
| Anthracene, 2-met...  | 11.83 | 4.0     | ng    | 203438   | 4                     | 11.33 | 1021870 | 20.0 |
| Phenanthrene, 1-m...  | 11.86 | 5.1     | ng    | 258751   | 4                     | 11.33 | 1021870 | 20.0 |
| Phenanthrene, 2-m...  | 11.90 | 2.7     | ng    | 139852   | 4                     | 11.33 | 1021870 | 20.0 |
| 4H-Cyclopenta[def...  | 11.95 | 8.8     | ng    | 448260   | 4                     | 11.33 | 1021870 | 20.0 |
| 1H-Cyclopropa[l]p...  | 11.96 | 2.8     | ng    | 143141   | 4                     | 11.33 | 1021870 | 20.0 |
| Tricyclo[8.2.2.2(...  | 12.13 | 2.9     | ng    | 149727   | 4                     | 11.33 | 1021870 | 20.0 |
| 9,10-Dimethylanthe... | 12.38 | 3.9     | ng    | 200313   | 4                     | 11.33 | 1021870 | 20.0 |
| unknown12.42          | 12.42 | 7.7     | ng    | 392256   | 4                     | 11.33 | 1021870 | 20.0 |
| Cyclopenta(def)ph...  | 12.47 | 2.5     | ng    | 126396   | 4                     | 11.33 | 1021870 | 20.0 |
| 11H-Benzo[a]fluorene  | 13.10 | 3.6     | ng    | 353043   | 5                     | 13.97 | 1943220 | 20.0 |
| 11H-Benzo[b]fluorene  | 13.16 | 2.5     | ng    | 241161   | 5                     | 13.97 | 1943220 | 20.0 |