

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
 Data File : BF101583.D  
 Acq On : 20 Dec 2017 22:14  
 Operator : SJ/JU  
 Sample : I6981-04 2X  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PHMHA-WCD

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-BF121417.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         | 5.004       | 493           | 496         | 515          | rBV      | 159128         | 232910        | 12.08%          | 0.738%        |
| 2         | 5.410       | 562           | 565         | 592          | rBV      | 1400563        | 1888351       | 97.97%          | 5.982%        |
| 3         | 6.434       | 736           | 739         | 758          | rBV      | 1587734        | 1927465       | 100.00%         | 6.106%        |
| 4         | 6.563       | 758           | 761         | 768          | rVV      | 1799350        | 1829950       | 94.94%          | 5.797%        |
| 5         | 6.616       | 768           | 770         | 777          | rVB3     | 22417          | 28233         | 1.46%           | 0.089%        |
| 6         | 6.792       | 797           | 800         | 816          | rBV      | 997471         | 938757        | 48.70%          | 2.974%        |
| 7         | 6.945       | 823           | 826         | 840          | rVB      | 1360332        | 1434392       | 74.42%          | 4.544%        |
| 8         | 7.363       | 893           | 897         | 915          | rBV      | 1107564        | 1196877       | 62.10%          | 3.792%        |
| 9         | 8.075       | 1015          | 1018        | 1030         | rBV      | 1200724        | 1217281       | 63.15%          | 3.856%        |
| 10        | 8.657       | 1114          | 1117        | 1122         | rBV      | 42549          | 34675         | 1.80%           | 0.110%        |
| 11        | 8.798       | 1138          | 1141        | 1145         | rBV      | 43092          | 48346         | 2.51%           | 0.153%        |
| 12        | 8.892       | 1154          | 1157        | 1160         | rBV      | 44332          | 37086         | 1.92%           | 0.117%        |
| 13        | 9.086       | 1187          | 1190        | 1197         | rBV      | 24988          | 21251         | 1.10%           | 0.067%        |
| 14        | 9.151       | 1197          | 1201        | 1209         | rVV      | 2014113        | 1862944       | 96.65%          | 5.902%        |
| 15        | 9.222       | 1209          | 1213        | 1216         | rVB      | 72826          | 66495         | 3.45%           | 0.211%        |
| 16        | 9.257       | 1216          | 1219        | 1223         | rBV2     | 33025          | 33570         | 1.74%           | 0.106%        |
| 17        | 9.351       | 1230          | 1235        | 1240         | rBV2     | 13886          | 19604         | 1.02%           | 0.062%        |
| 18        | 9.416       | 1243          | 1246        | 1251         | rBV2     | 37010          | 53202         | 2.76%           | 0.169%        |
| 19        | 9.492       | 1255          | 1259        | 1262         | rBV2     | 64427          | 73792         | 3.83%           | 0.234%        |
| 20        | 9.522       | 1262          | 1264        | 1265         | rVV      | 33399          | 29625         | 1.54%           | 0.094%        |
| 21        | 9.545       | 1265          | 1268        | 1275         | rVB2     | 85662          | 118626        | 6.15%           | 0.376%        |
| 22        | 9.610       | 1275          | 1279        | 1281         | rBV2     | 33174          | 37560         | 1.95%           | 0.119%        |
| 23        | 9.698       | 1290          | 1294        | 1298         | rBV2     | 68679          | 78244         | 4.06%           | 0.248%        |
| 24        | 9.751       | 1300          | 1303        | 1305         | rVV      | 89031          | 71544         | 3.71%           | 0.227%        |
| 25        | 9.833       | 1305          | 1317        | 1320         | rVV      | 1179776        | 1130657       | 58.66%          | 3.582%        |
| 26        | 9.863       | 1320          | 1322        | 1329         | rVB      | 218328         | 205933        | 10.68%          | 0.652%        |
| 27        | 9.933       | 1330          | 1334        | 1338         | rBV3     | 23891          | 31566         | 1.64%           | 0.100%        |
| 28        | 9.986       | 1339          | 1343        | 1345         | rBV2     | 25998          | 29456         | 1.53%           | 0.093%        |
| 29        | 10.039      | 1348          | 1352        | 1355         | rVV2     | 111084         | 112715        | 5.85%           | 0.357%        |
| 30        | 10.075      | 1355          | 1358        | 1361         | rVB3     | 45118          | 46597         | 2.42%           | 0.148%        |
| 31        | 10.145      | 1366          | 1370        | 1373         | rBV      | 36042          | 42946         | 2.23%           | 0.136%        |
| 32        | 10.180      | 1373          | 1376        | 1381         | rVB3     | 33633          | 41097         | 2.13%           | 0.130%        |
| 33        | 10.245      | 1381          | 1387        | 1391         | rBV2     | 119811         | 152356        | 7.90%           | 0.483%        |
| 34        | 10.380      | 1407          | 1410        | 1416         | rVB      | 146700         | 149955        | 7.78%           | 0.475%        |

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Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
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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:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 10.469 | 1422 | 1425 | 1427 | rVB  | 67964   | 63075   | 3.27%  | 0.200% |
| 36 | 10.516 | 1431 | 1433 | 1435 | rBV3 | 24220   | 22476   | 1.17%  | 0.071% |
| 37 | 10.539 | 1435 | 1437 | 1441 | rVB2 | 47937   | 45390   | 2.35%  | 0.144% |
| 38 | 10.628 | 1448 | 1452 | 1465 | rBV  | 749176  | 833232  | 43.23% | 2.640% |
| 39 | 10.722 | 1465 | 1468 | 1473 | rBV  | 141996  | 188034  | 9.76%  | 0.596% |
| 40 | 10.816 | 1481 | 1484 | 1487 | rBV5 | 19460   | 19753   | 1.02%  | 0.063% |
| 41 | 10.933 | 1495 | 1504 | 1506 | rBV2 | 57134   | 103887  | 5.39%  | 0.329% |
| 42 | 11.057 | 1520 | 1525 | 1527 | rBV4 | 33446   | 54460   | 2.83%  | 0.173% |
| 43 | 11.080 | 1527 | 1529 | 1535 | rVV4 | 32304   | 45976   | 2.39%  | 0.146% |
| 44 | 11.139 | 1536 | 1539 | 1541 | rVV  | 80207   | 70624   | 3.66%  | 0.224% |
| 45 | 11.169 | 1541 | 1544 | 1547 | rVV  | 123699  | 119558  | 6.20%  | 0.379% |
| 46 | 11.198 | 1547 | 1549 | 1551 | rVV2 | 61314   | 57912   | 3.00%  | 0.183% |
| 47 | 11.222 | 1551 | 1553 | 1557 | rVB  | 89846   | 79082   | 4.10%  | 0.251% |
| 48 | 11.322 | 1566 | 1570 | 1572 | rBV  | 1020615 | 908091  | 47.11% | 2.877% |
| 49 | 11.345 | 1572 | 1574 | 1579 | rVV  | 1471854 | 1234772 | 64.06% | 3.912% |
| 50 | 11.398 | 1580 | 1583 | 1588 | rVB  | 369827  | 356512  | 18.50% | 1.129% |
| 51 | 11.563 | 1605 | 1611 | 1614 | rBV2 | 146419  | 178593  | 9.27%  | 0.566% |
| 52 | 11.592 | 1614 | 1616 | 1618 | rVV  | 61373   | 46077   | 2.39%  | 0.146% |
| 53 | 11.822 | 1652 | 1655 | 1658 | rBV  | 170994  | 163848  | 8.50%  | 0.519% |
| 54 | 11.851 | 1658 | 1660 | 1665 | rVV  | 202209  | 199028  | 10.33% | 0.631% |
| 55 | 11.904 | 1665 | 1669 | 1672 | rVV  | 93507   | 96317   | 5.00%  | 0.305% |
| 56 | 11.939 | 1672 | 1675 | 1677 | rVV  | 232796  | 260160  | 13.50% | 0.824% |
| 57 | 11.957 | 1677 | 1678 | 1683 | rVB  | 115217  | 104997  | 5.45%  | 0.333% |
| 58 | 11.998 | 1683 | 1685 | 1690 | rBV2 | 61434   | 75901   | 3.94%  | 0.240% |
| 59 | 12.116 | 1702 | 1705 | 1709 | rBV  | 129433  | 113773  | 5.90%  | 0.360% |
| 60 | 12.163 | 1710 | 1713 | 1716 | rBV2 | 73695   | 73961   | 3.84%  | 0.234% |
| 61 | 12.269 | 1725 | 1731 | 1734 | rBV3 | 49230   | 102067  | 5.30%  | 0.323% |
| 62 | 12.298 | 1734 | 1736 | 1738 | rVB  | 38606   | 31197   | 1.62%  | 0.099% |
| 63 | 12.374 | 1746 | 1749 | 1750 | rBV  | 113346  | 104563  | 5.42%  | 0.331% |
| 64 | 12.474 | 1761 | 1766 | 1769 | rBV2 | 74027   | 108496  | 5.63%  | 0.344% |
| 65 | 12.539 | 1773 | 1777 | 1781 | rBV  | 1635865 | 1482478 | 76.91% | 4.697% |
| 66 | 12.604 | 1786 | 1788 | 1792 | rBV2 | 35582   | 53677   | 2.78%  | 0.170% |
| 67 | 12.692 | 1800 | 1803 | 1805 | rBV2 | 88044   | 77358   | 4.01%  | 0.245% |
| 68 | 12.769 | 1810 | 1816 | 1822 | rBV  | 1378989 | 1609990 | 83.53% | 5.101% |
| 69 | 12.821 | 1822 | 1825 | 1829 | rVB  | 82587   | 87161   | 4.52%  | 0.276% |
| 70 | 12.904 | 1835 | 1839 | 1842 | rBV  | 1336581 | 1190283 | 61.75% | 3.771% |
| 71 | 12.939 | 1843 | 1845 | 1848 | rVB  | 77357   | 70099   | 3.64%  | 0.222% |

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ClientSampleId :  
PHMHA-WCD

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

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

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 72 | 12.992 | 1849 | 1854 | 1856 | rBV2 | 89967   | 148033  | 7.68%  | 0.469% |
| 73 | 13.074 | 1865 | 1868 | 1869 | rBV2 | 81999   | 81670   | 4.24%  | 0.259% |
| 74 | 13.098 | 1869 | 1872 | 1877 | rVB2 | 208571  | 240196  | 12.46% | 0.761% |
| 75 | 13.163 | 1881 | 1883 | 1885 | rVV  | 157715  | 119742  | 6.21%  | 0.379% |
| 76 | 13.204 | 1888 | 1890 | 1894 | rVB  | 108871  | 109610  | 5.69%  | 0.347% |
| 77 | 13.298 | 1903 | 1906 | 1908 | rBV  | 79605   | 72424   | 3.76%  | 0.229% |
| 78 | 13.327 | 1909 | 1911 | 1917 | rVB2 | 89574   | 99552   | 5.16%  | 0.315% |
| 79 | 13.616 | 1958 | 1960 | 1963 | rVB  | 97975   | 90127   | 4.68%  | 0.286% |
| 80 | 13.721 | 1976 | 1978 | 1980 | rBV2 | 128249  | 118662  | 6.16%  | 0.376% |
| 81 | 13.792 | 1985 | 1990 | 1994 | rVV3 | 199476  | 369409  | 19.17% | 1.170% |
| 82 | 13.827 | 1994 | 1996 | 2002 | rVB  | 135271  | 147080  | 7.63%  | 0.466% |
| 83 | 13.921 | 2010 | 2012 | 2014 | rBV  | 99900   | 80508   | 4.18%  | 0.255% |
| 84 | 13.968 | 2015 | 2020 | 2023 | rVV2 | 1090843 | 1651982 | 85.71% | 5.234% |
| 85 | 13.998 | 2023 | 2025 | 2029 | rVB  | 702918  | 583222  | 30.26% | 1.848% |
| 86 | 14.074 | 2036 | 2038 | 2041 | rBV  | 111824  | 106278  | 5.51%  | 0.337% |
| 87 | 15.015 | 2194 | 2198 | 2200 | rBV  | 496426  | 673011  | 34.92% | 2.132% |
| 88 | 15.315 | 2246 | 2249 | 2252 | rBV  | 289543  | 307443  | 15.95% | 0.974% |
| 89 | 15.380 | 2257 | 2260 | 2267 | rVB  | 362240  | 474578  | 24.62% | 1.504% |
| 90 | 15.439 | 2267 | 2270 | 2273 | rBV  | 418335  | 434379  | 22.54% | 1.376% |

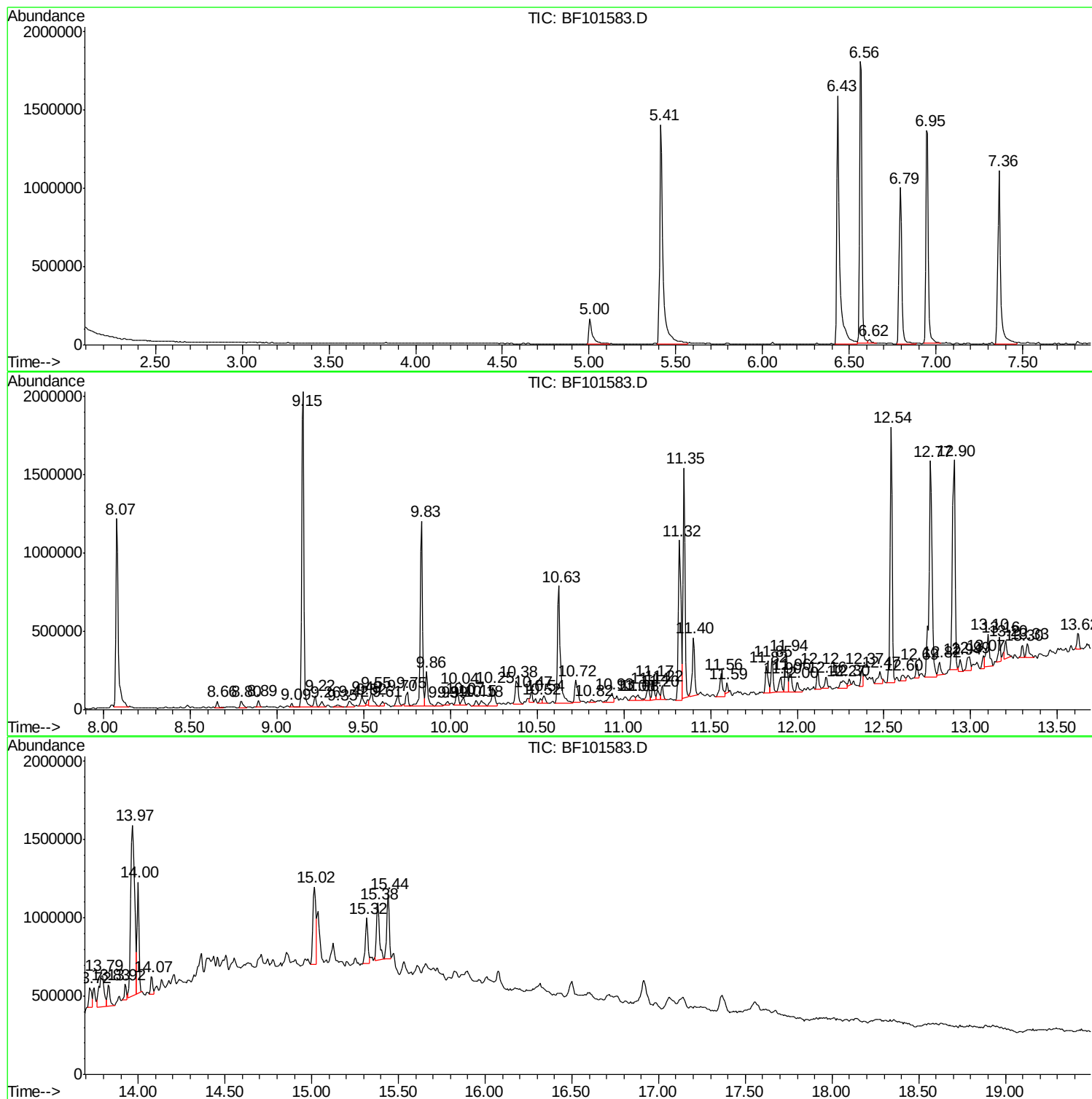
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ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.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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Misc :  
ALS Vial : 25 Sample Multiplier: 1

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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

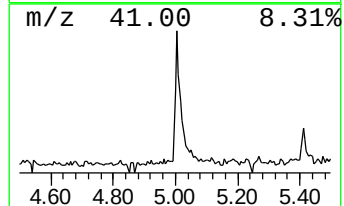
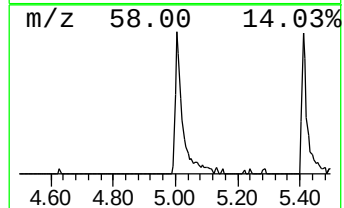
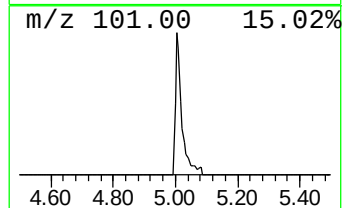
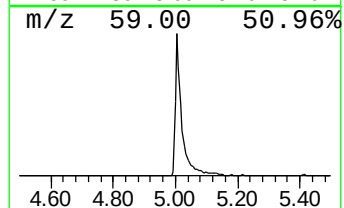
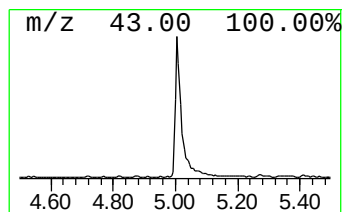
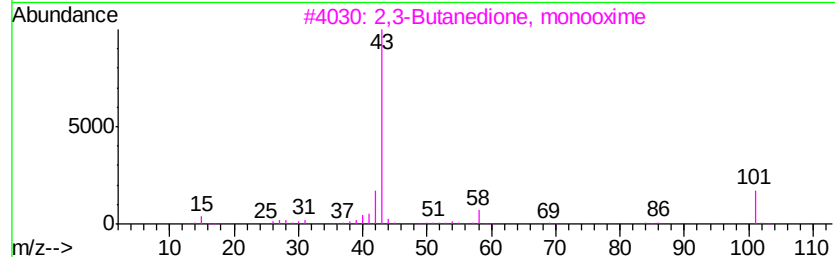
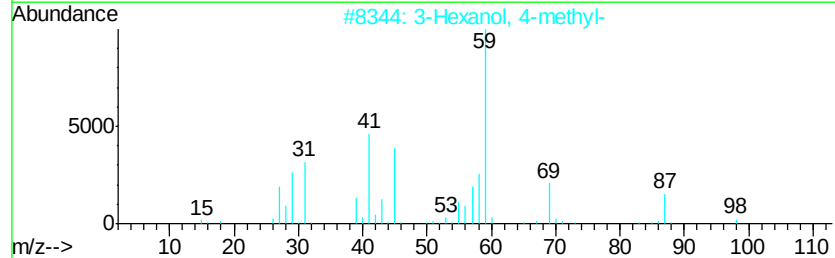
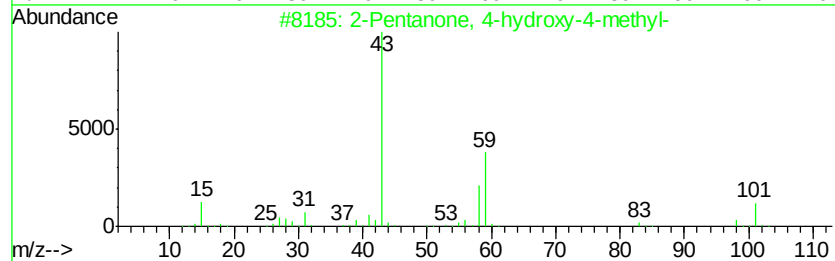
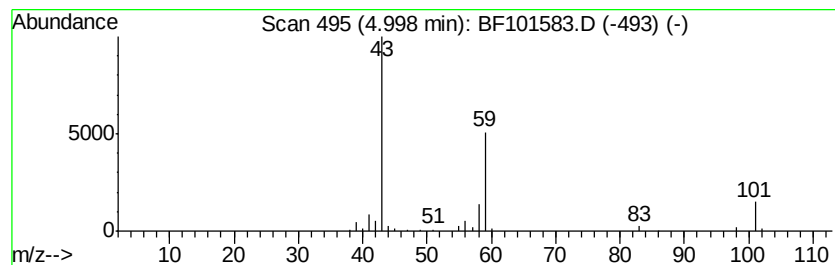
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.00 | 4.96 ng | 232910 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2 | 64   |
| 2    |    |   | 3-Hexanol, 4-methyl-             | 116 | C7H16O  | 000615-29-2 | 33   |
| 3    |    |   | 2,3-Butanedione, monooxime       | 101 | C4H7NO2 | 000057-71-6 | 9    |
| 4    |    |   | Propane, 1-ethoxy-2-methyl-      | 102 | C6H14O  | 000627-02-1 | 9    |
| 5    |    |   | 2-Propanone, 1-(1-methylethoxy)- | 116 | C6H12O2 | 042781-12-4 | 9    |



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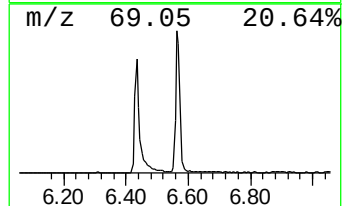
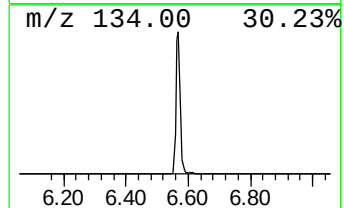
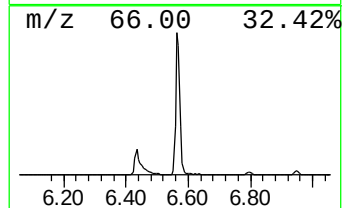
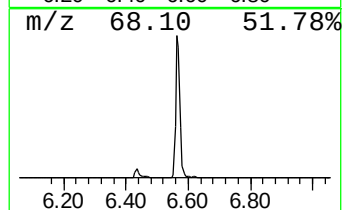
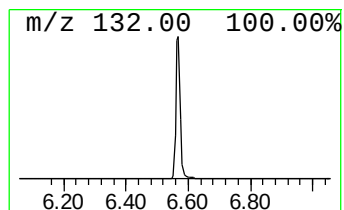
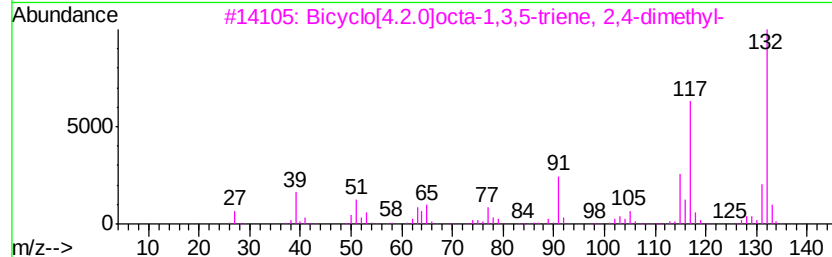
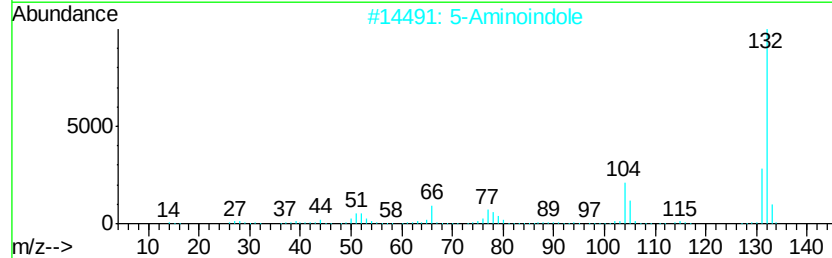
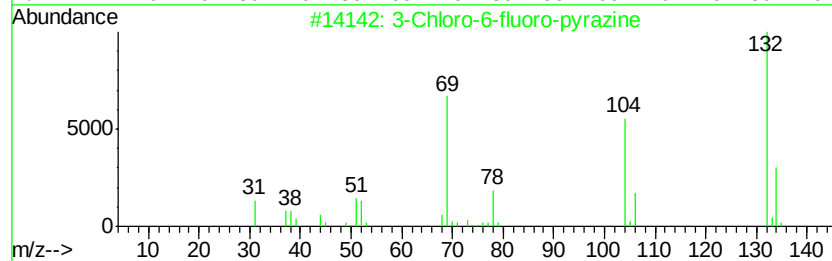
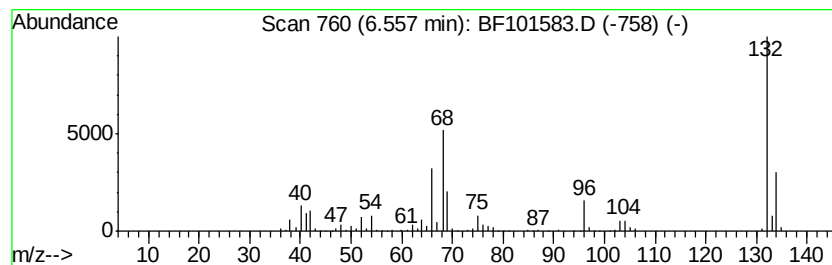
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 unknown6.56 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.56 | 38.99 ng | 1829950 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 35   |
| 2    |    | 5-Aminoindole                       | 132 | C8H8N2    | 005192-03-0  | 12   |
| 3    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 4    |    | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2    | 022291-85-6  | 9    |
| 5    |    | 1-Methylpyrrolo[1,2-a]pyrazine      | 132 | C8H8N2    | 064608-59-9  | 9    |



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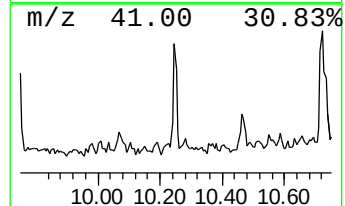
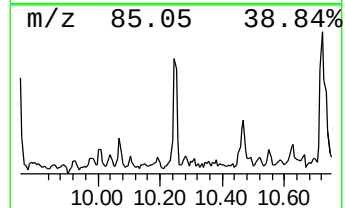
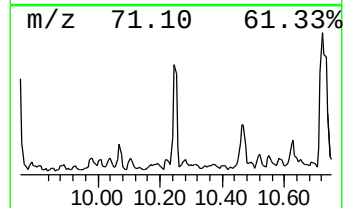
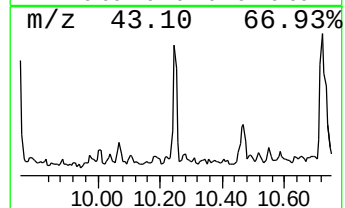
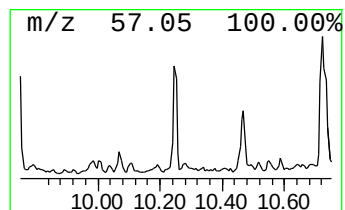
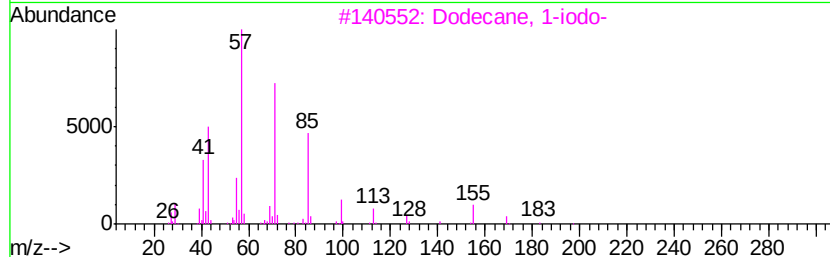
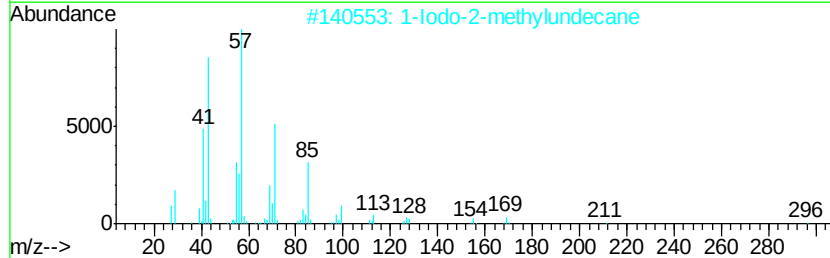
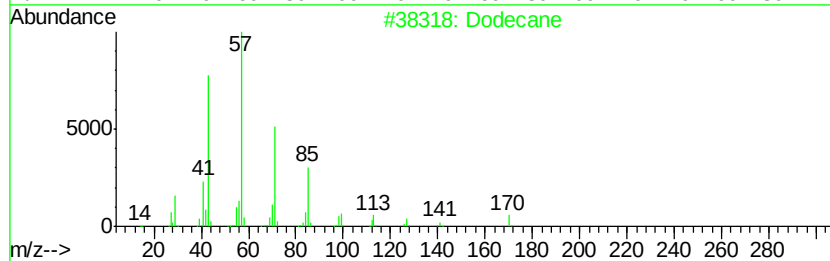
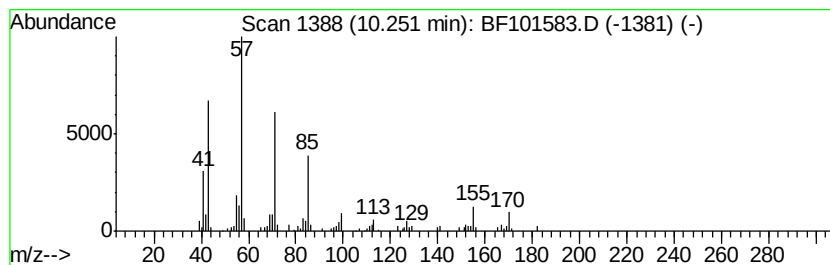
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Dodecane Concentration Rank 11

| R.T.    | EstConc                 | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------|--------------|------------------|----------------|
| 10.25   | 2.70 ng                 | 152356       | Acenaphthene-d10 | 9.83           |
| Hit# of | 5                       | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Dodecane                |              | 170 C12H26       | 000112-40-3 94 |
| 2       | 1-Iodo-2-methylundecane |              | 296 C12H25I      | 073105-67-6 72 |
| 3       | Dodecane, 1-iodo-       |              | 296 C12H25I      | 004292-19-7 72 |
| 4       | Hexadecane              |              | 226 C16H34       | 000544-76-3 68 |
| 5       | Pentadecane             |              | 212 C15H32       | 000629-62-9 64 |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

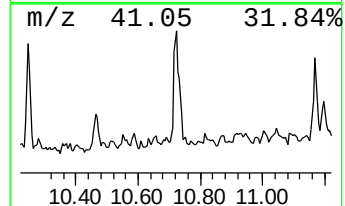
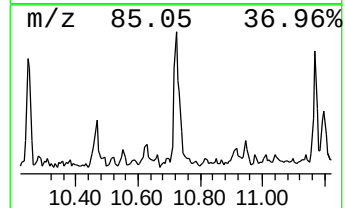
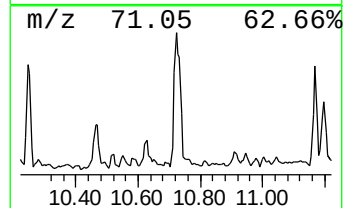
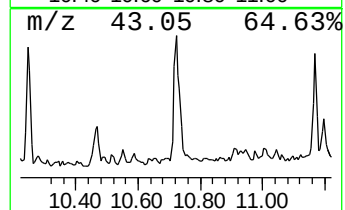
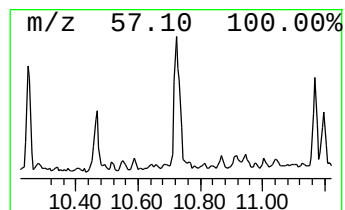
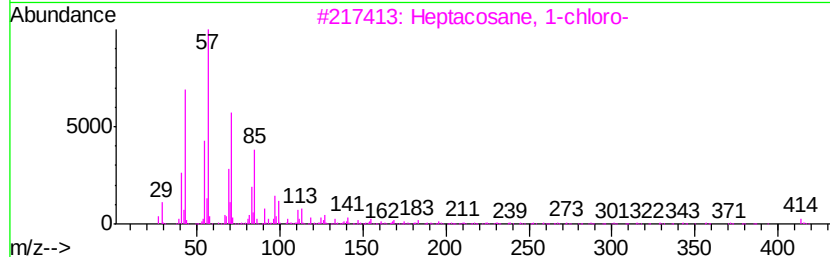
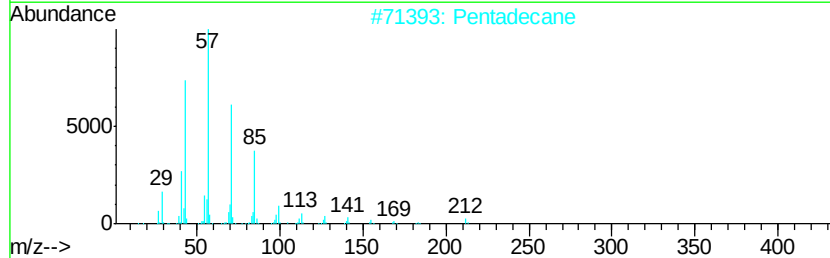
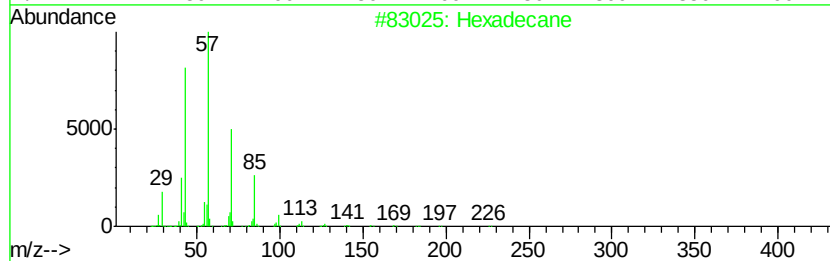
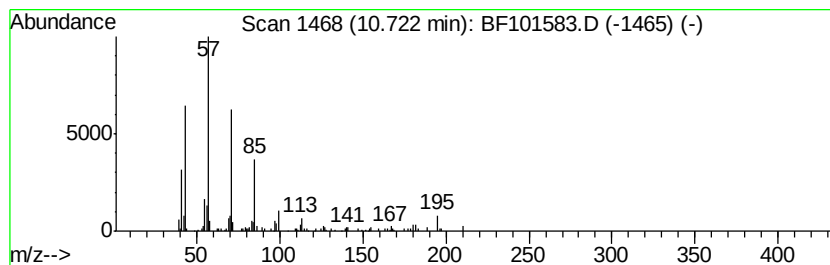
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Hexadecane Concentration Rank 8

| R.T.      | EstConc                | Area   | Relative to ISTD | R.T.        |      |
|-----------|------------------------|--------|------------------|-------------|------|
| 10.72     | 4.14 ng                | 188034 | Phenanthrene-d10 | 11.32       |      |
| Hit# of 5 | Tentative ID           | MW     | MolForm          | CAS#        | Qual |
| 1         | Hexadecane             | 226    | C16H34           | 000544-76-3 | 80   |
| 2         | Pentadecane            | 212    | C15H32           | 000629-62-9 | 80   |
| 3         | Heptacosane, 1-chloro- | 414    | C27H55Cl         | 062016-79-9 | 78   |
| 4         | Heptacosane            | 380    | C27H56           | 000593-49-7 | 78   |
| 5         | Tetradecane            | 198    | C14H30           | 000629-59-4 | 64   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

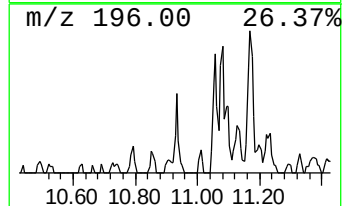
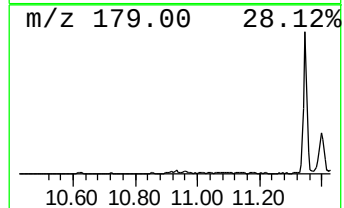
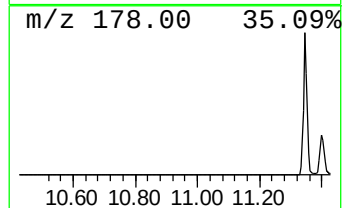
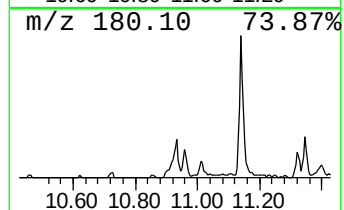
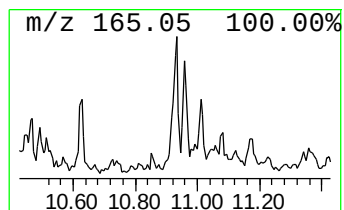
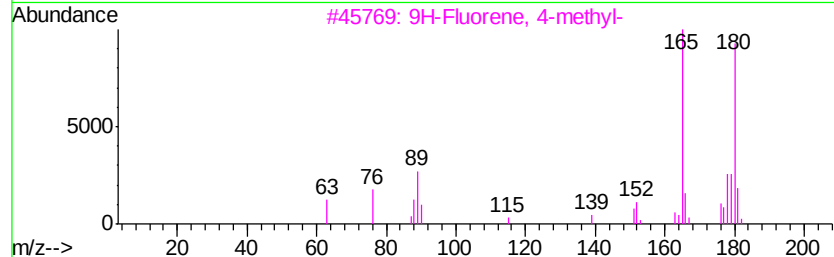
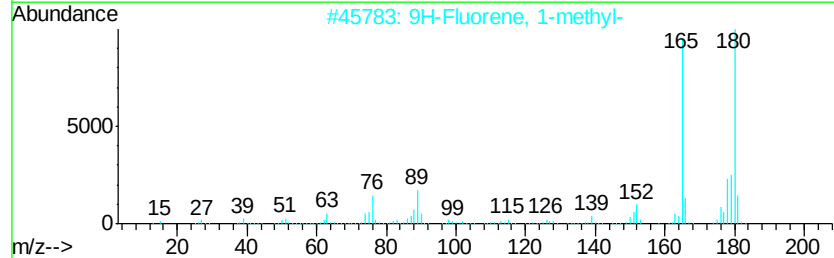
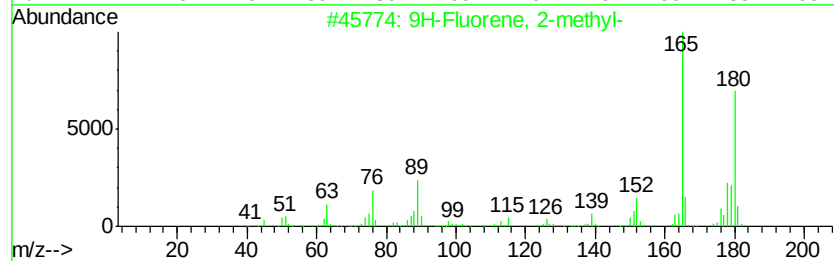
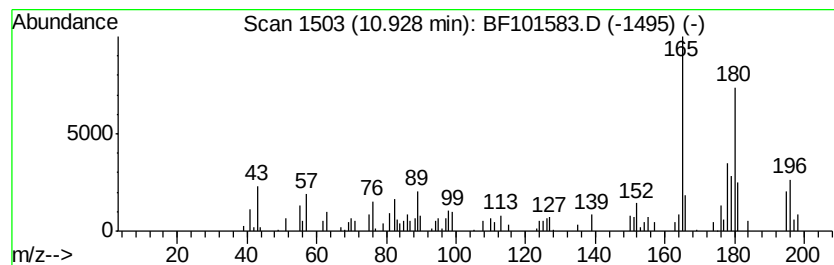
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 9H-Fluorene, 2-methyl- Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.93 | 2.29 ng | 103887 | Phenanthrene-d10 | 11.32 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | 9H-Fluorene, 2-methyl- | 180 | C14H12  | 001430-97-3 | 78   |
| 2    |    | 9H-Fluorene, 1-methyl- | 180 | C14H12  | 001730-37-6 | 78   |
| 3    |    | 9H-Fluorene, 4-methyl- | 180 | C14H12  | 001556-99-6 | 78   |
| 4    |    | 9H-Fluorene, 9-methyl- | 180 | C14H12  | 002523-37-7 | 60   |
| 5    |    | 9H-Fluorene, 3-methyl- | 180 | C14H12  | 002523-39-9 | 60   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

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

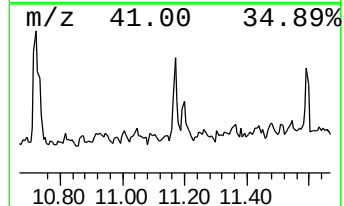
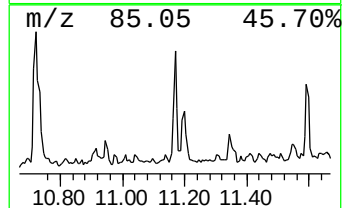
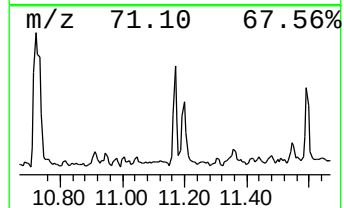
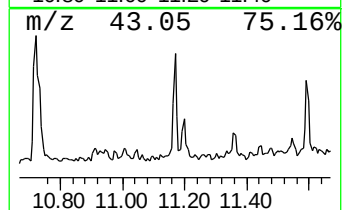
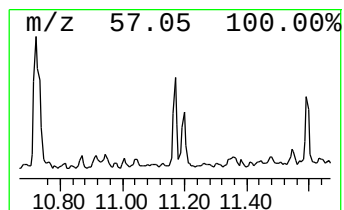
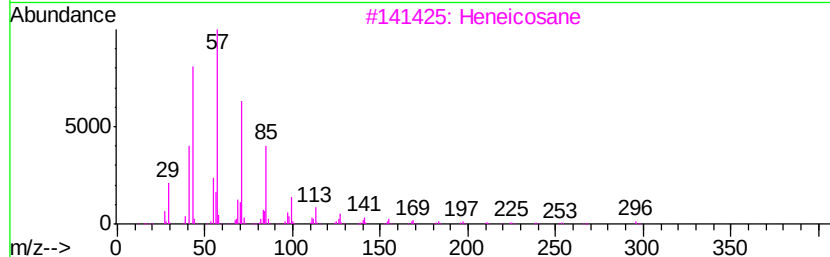
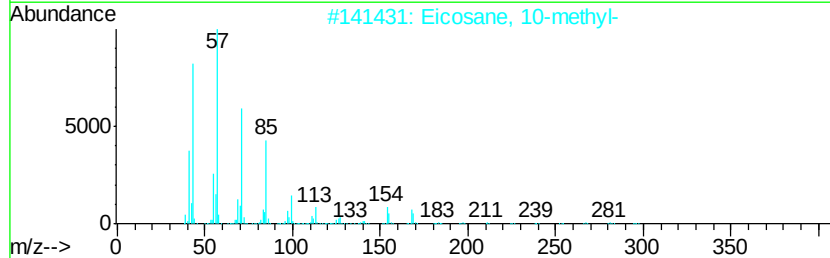
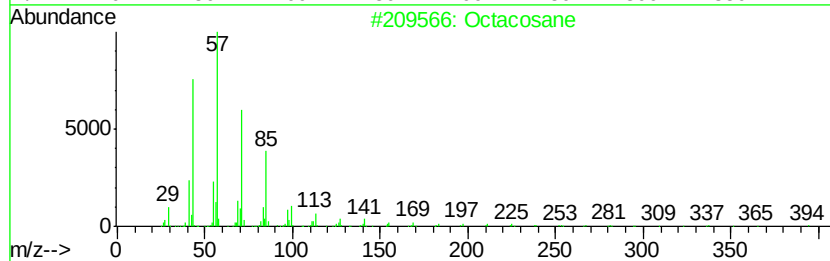
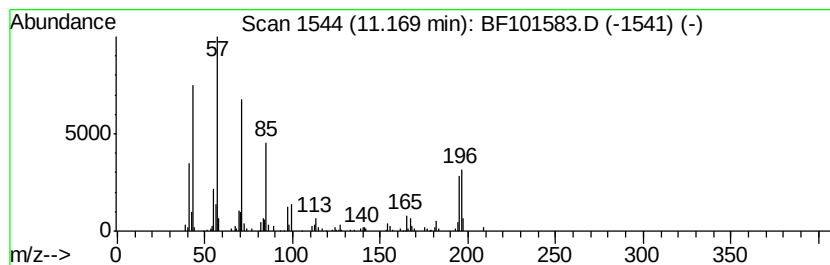
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 Octacosane Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.17 | 2.63 ng | 119558 | Phenanthrene-d10 | 11.32 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Octacosane                          | 394 | C28H58    | 000630-02-4  | 58   |
| 2    |    | Eicosane, 10-methyl-                | 296 | C21H44    | 054833-23-7  | 52   |
| 3    |    | Heneicosane                         | 296 | C21H44    | 000629-94-7  | 52   |
| 4    |    | Dodecane, 2-methyl-6-propyl-        | 226 | C16H34    | 055045-08-4  | 52   |
| 5    |    | Sulfurous acid, 2-propyl tridecy... | 306 | C16H34O3S | 1000309-12-4 | 52   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
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Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

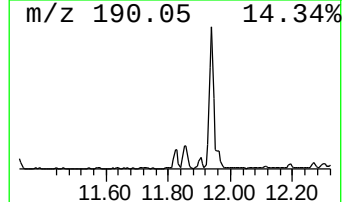
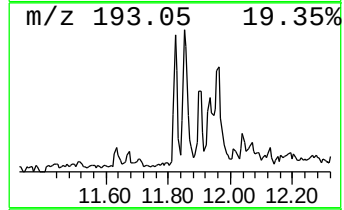
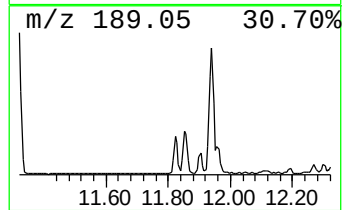
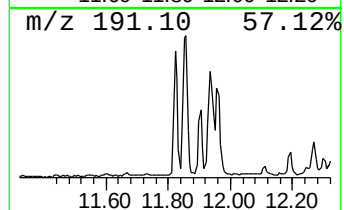
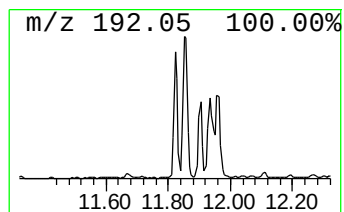
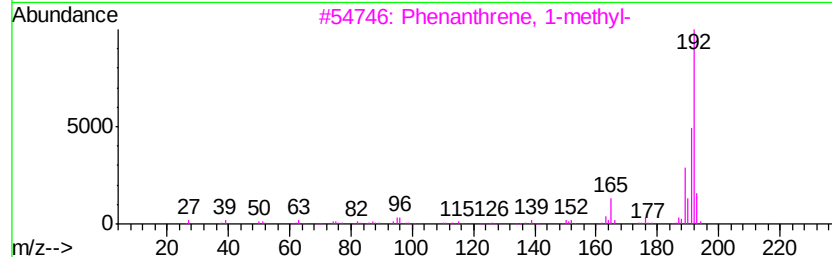
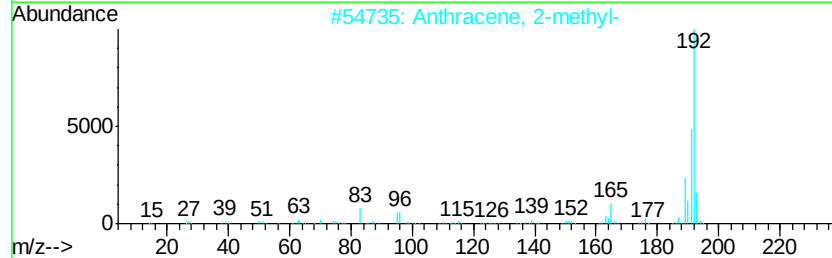
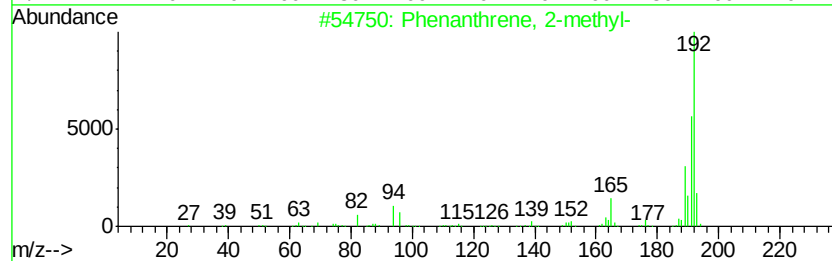
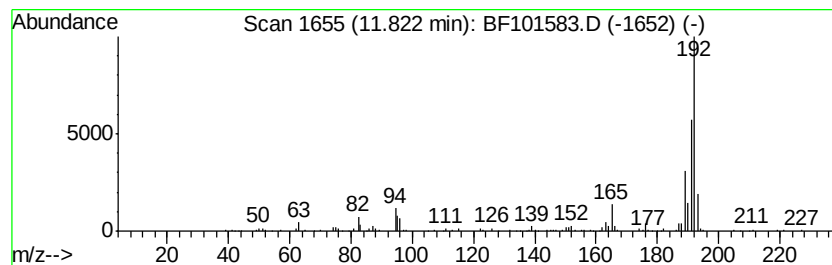
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 9

| R.T.      | EstConc                             | Area   | Relative to ISTD |             | R.T.  |
|-----------|-------------------------------------|--------|------------------|-------------|-------|
| 11.82     | 3.61 ng                             | 163848 | Phenanthrene-d10 |             | 11.32 |
| 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 | 97    |
| 3         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 97    |
| 4         | 1H-Cyclopropa[l]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 96    |
| 5         | 1H-Indene, 2-phenyl-                | 192    | C15H12           | 004505-48-0 | 94    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

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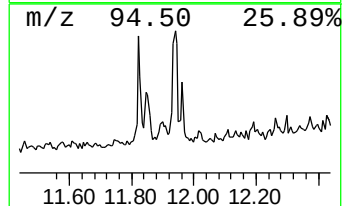
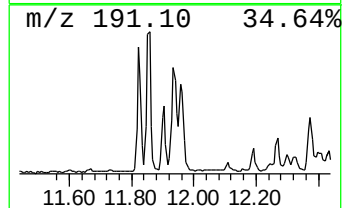
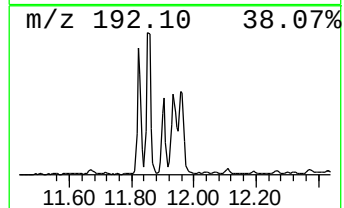
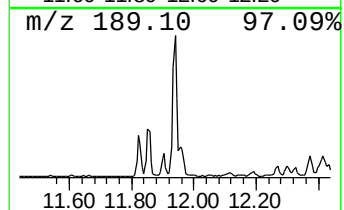
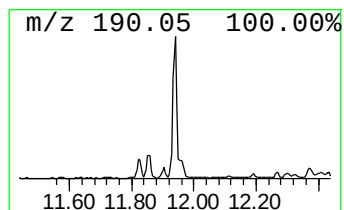
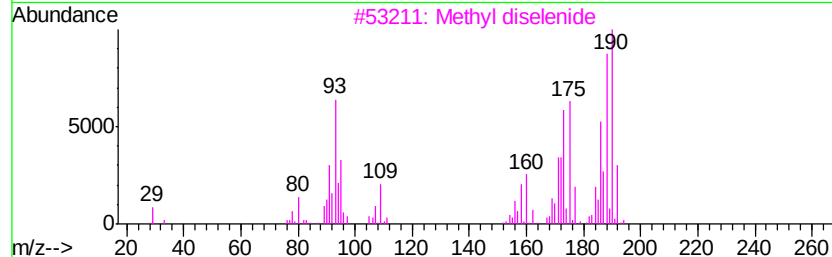
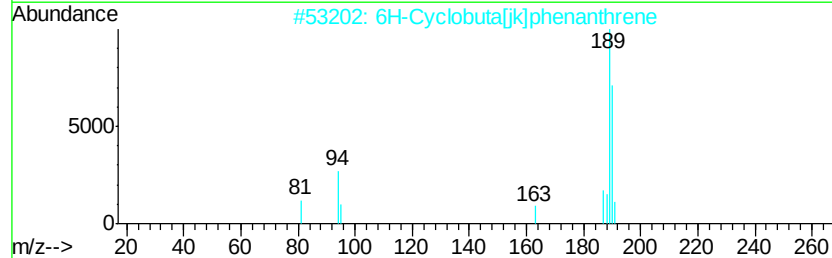
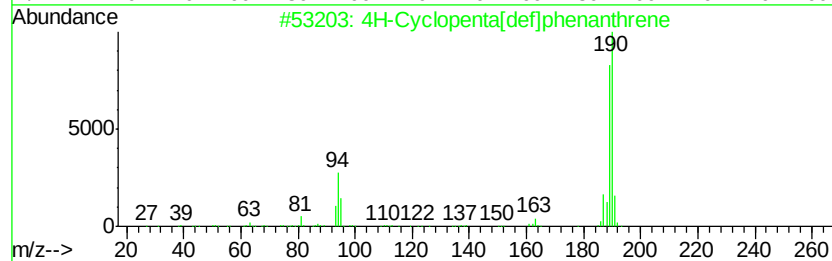
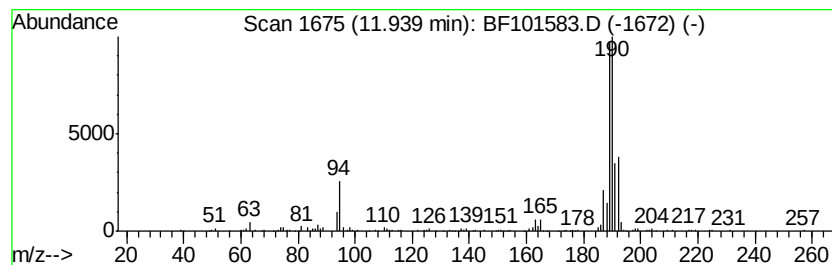
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TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.94 | 5.73 ng | 260160 | Phenanthrene-d10 | 11.32 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10       | 000203-64-5  | 81   |
| 2    |    | 6H-Cyclobuta[jk]phenanthrene        | 190 | C15H10       | 083469-43-6  | 59   |
| 3    |    | Methyl diselenide                   | 190 | C2H6Se2      | 007101-31-7  | 46   |
| 4    |    | 2,2'-Bis(4,5-dimethylimidazole)     | 190 | C10H14N4     | 069286-06-2  | 42   |
| 5    |    | 1-Aminocyclopentanecarboxylic ac... | 297 | C11H17Cl2N04 | 1000329-17-6 | 17   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

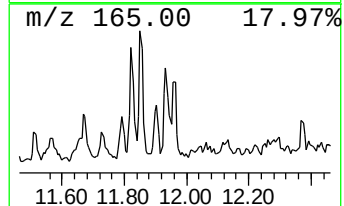
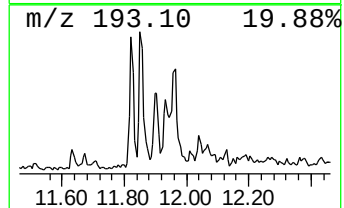
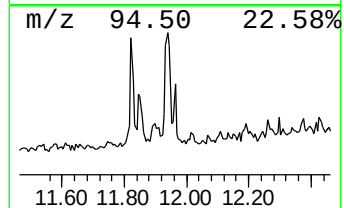
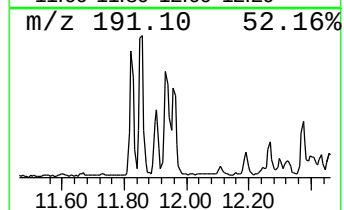
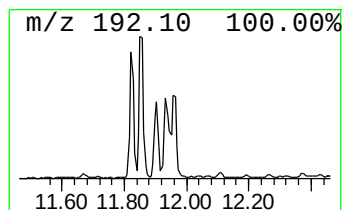
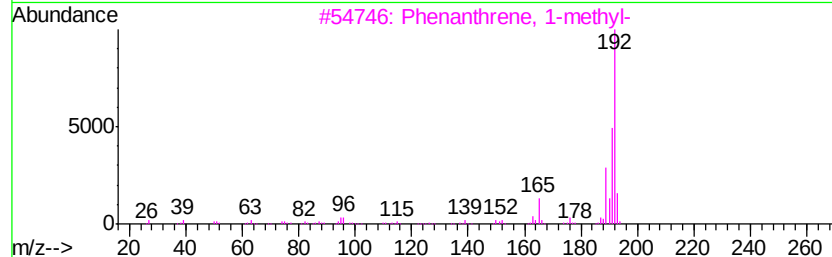
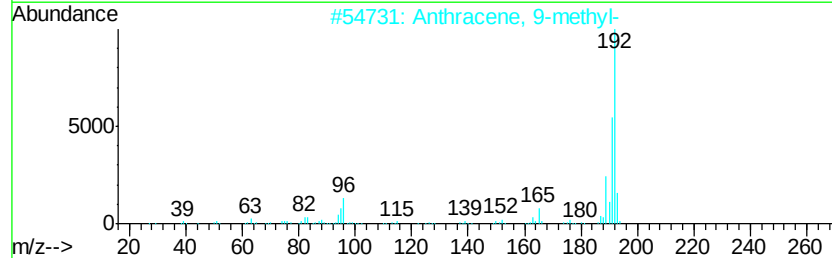
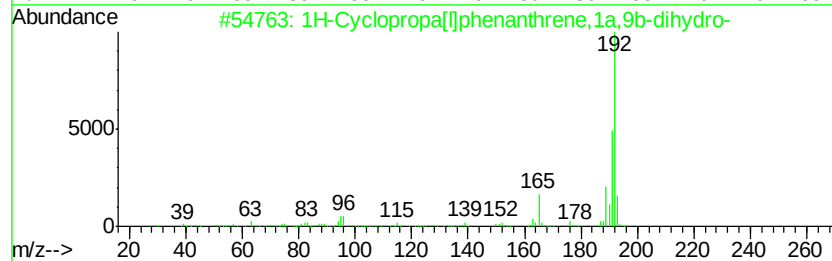
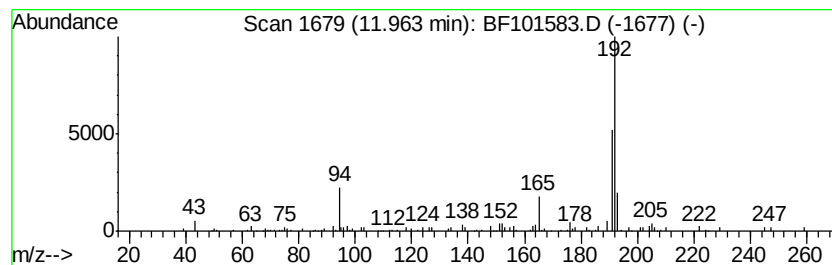
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 1H-Cyclopropa[l]phenanthren... Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.96 | 2.31 ng | 104997 | Phenanthrene-d10 | 11.32 |

| Hit# | of | 5 | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Cyclopropa[l]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 83   |
| 2    |    |   | Anthracene, 9-methyl-                 | 192 | C15H12  | 000779-02-2 | 83   |
| 3    |    |   | Phenanthrene, 1-methyl-               | 192 | C15H12  | 000832-69-9 | 83   |
| 4    |    |   | Phenanthrene, 2-methyl-               | 192 | C15H12  | 002531-84-2 | 83   |
| 5    |    |   | Phenanthrene, 4-methyl-               | 192 | C15H12  | 000832-64-4 | 83   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

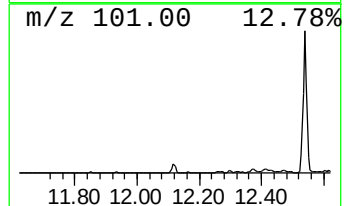
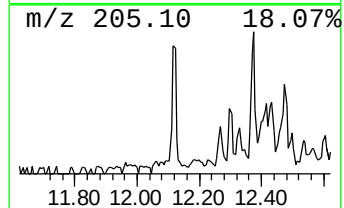
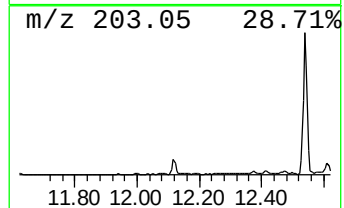
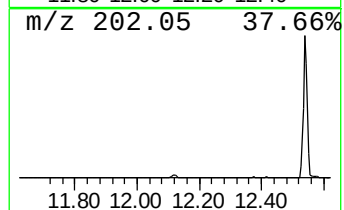
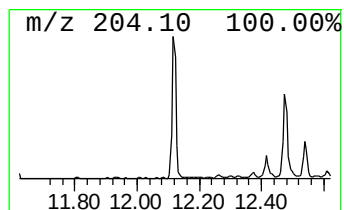
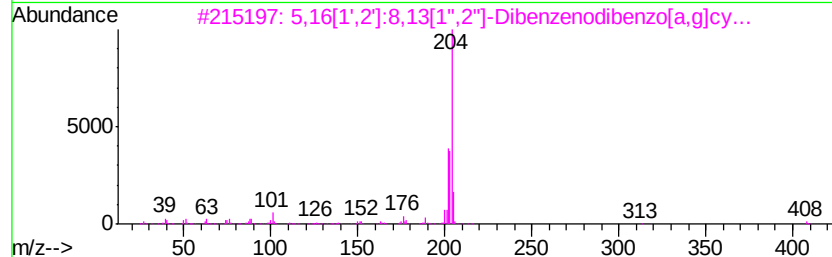
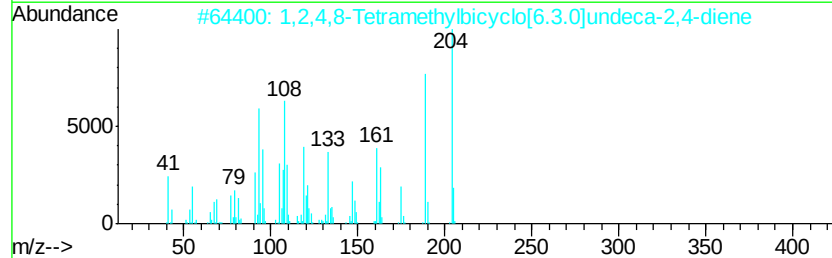
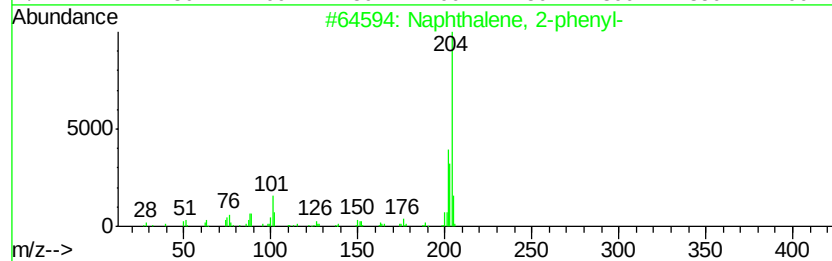
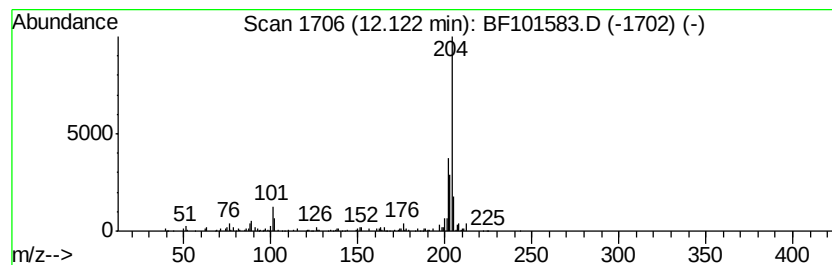
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Naphthalene, 2-phenyl- Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.12 | 2.51 ng | 113773 | Phenanthrene-d10 | 11.32 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-phenyl-               | 204 | C16H12  | 000612-94-2 | 93   |
| 2    |    | 1,2,4,8-Tetramethylbicyclo[6.3.0...] | 204 | C15H24  | 137235-51-9 | 90   |
| 3    |    | 5,16[1',2']:8,13[1'',2'']-Dibenz...  | 408 | C32H24  | 005672-97-9 | 86   |
| 4    |    | Naphthalene, 1-phenyl-               | 204 | C16H12  | 000605-02-7 | 83   |
| 5    |    | Tricyclo[8.2.2.2(4,7)]hexadeca-2...  | 204 | C16H12  | 006572-60-7 | 81   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

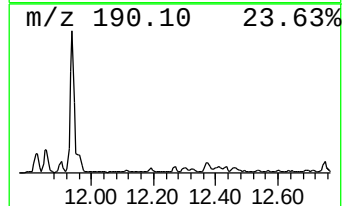
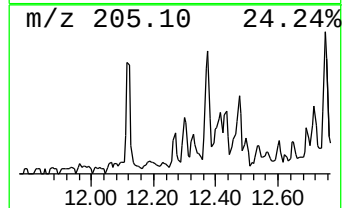
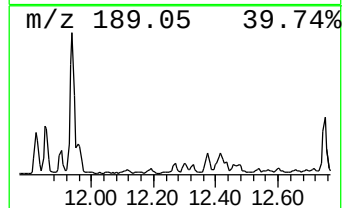
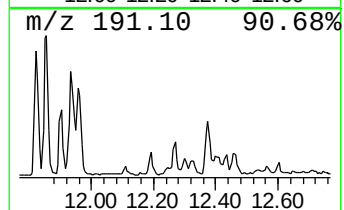
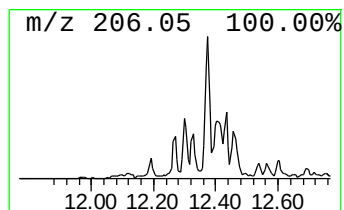
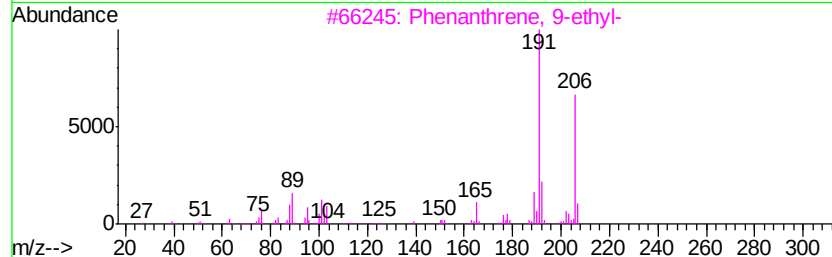
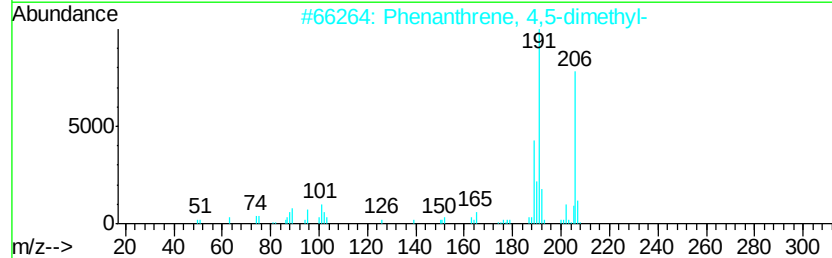
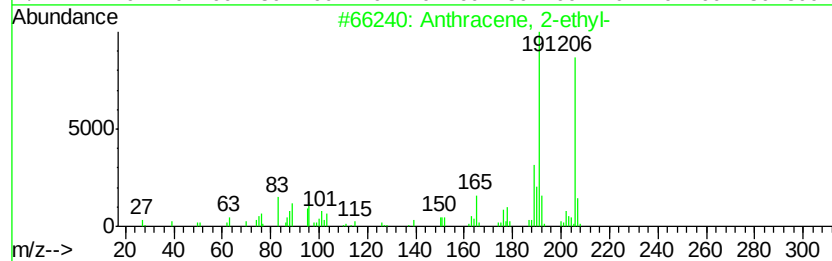
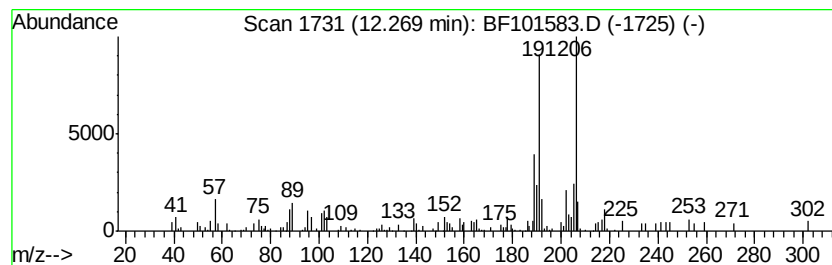
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Anthracene, 2-ethyl- Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.27 | 2.25 ng | 102067 | Phenanthrene-d10 | 11.32 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 2-ethyl-        | 206 | C16H14  | 052251-71-5 | 93   |
| 2    |    | Phenanthrene, 4,5-dimethyl- | 206 | C16H14  | 003674-69-9 | 93   |
| 3    |    | Phenanthrene, 9-ethyl-      | 206 | C16H14  | 003674-75-7 | 92   |
| 4    |    | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 87   |
| 5    |    | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 83   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

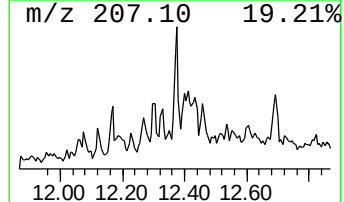
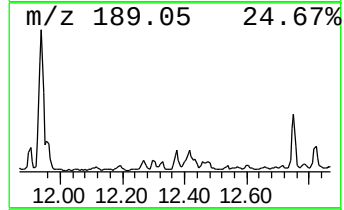
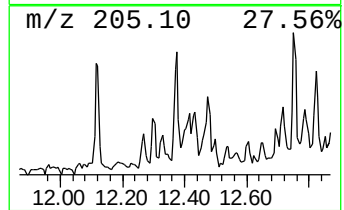
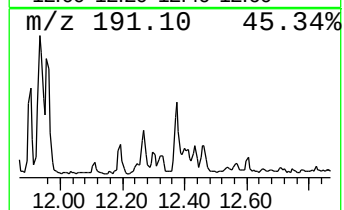
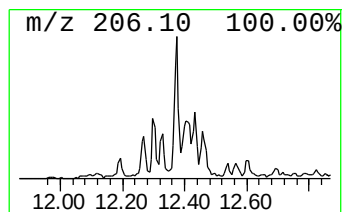
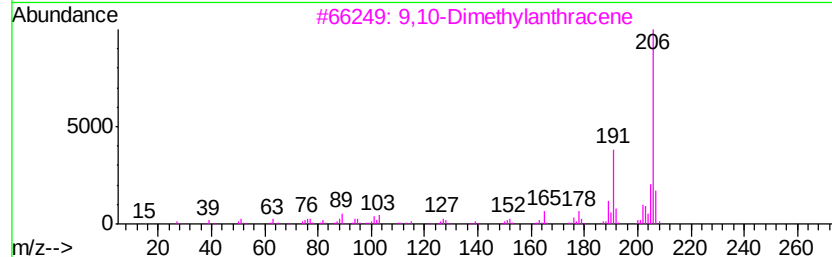
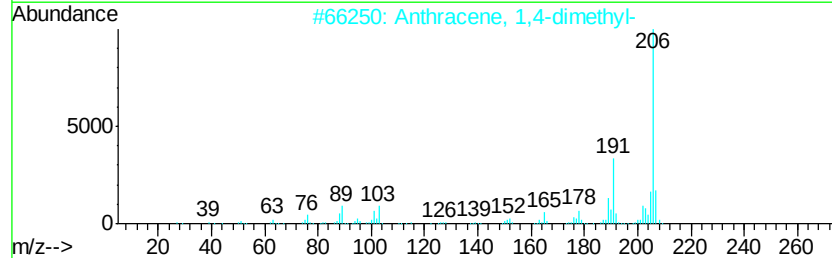
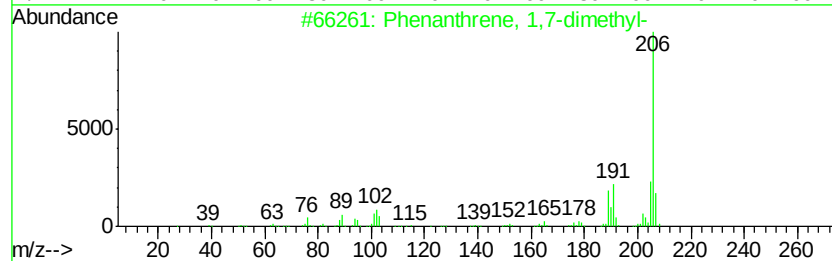
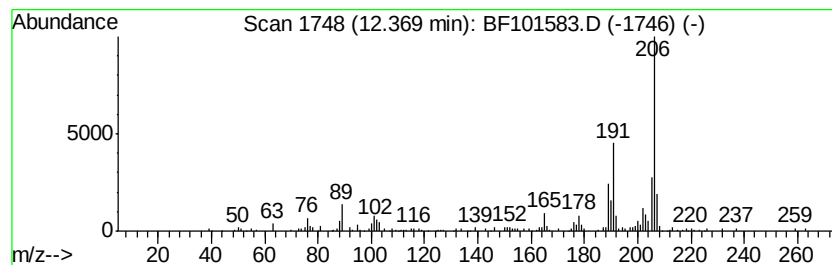
Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.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 Phenanthrene, 1,7-dimethyl- Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD            |     | R.T.    |             |      |
|-------|---------|--------|-----------------------------|-----|---------|-------------|------|
| 12.37 | 2.30 ng | 104563 | Phenanthrene-d10            |     | 11.32   |             |      |
| Hit#  | of      | 5      | Tentative ID                | MW  | MolForm | CAS#        | Qual |
| 1     |         |        | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 91   |
| 2     |         |        | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 91   |
| 3     |         |        | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 91   |
| 4     |         |        | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 89   |
| 5     |         |        | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 83   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

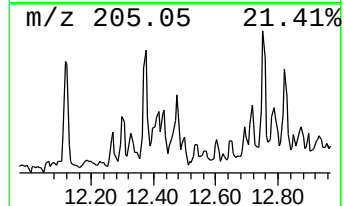
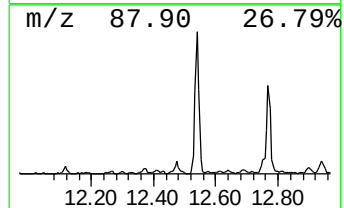
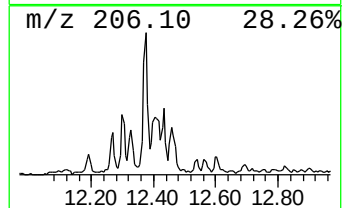
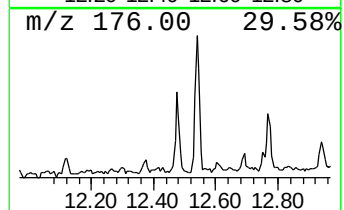
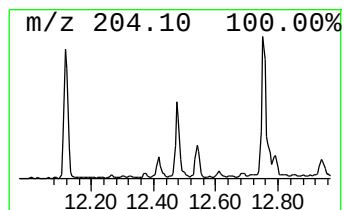
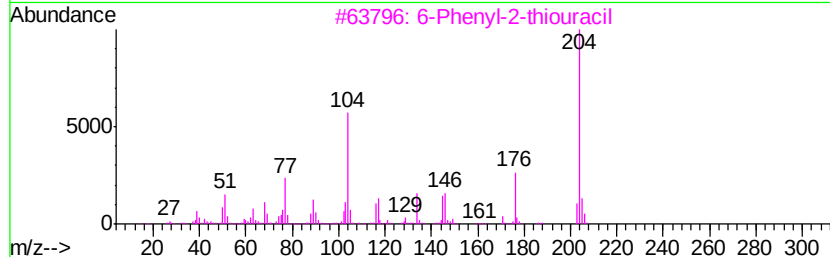
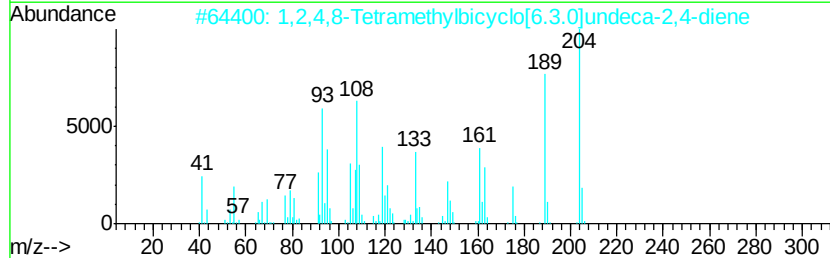
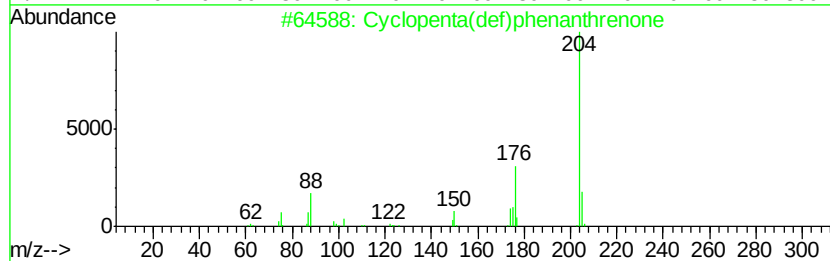
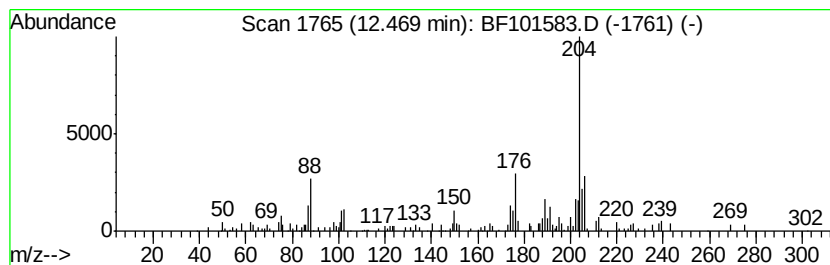
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Cyclopenta(def)phenanthrenone Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.47 | 2.39 ng | 108496 | Phenanthrene-d10 | 11.32 |

| Hit# | of | Tentative ID                         | MW  | MolForm   | CAS#        | Qual |
|------|----|--------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrenone        | 204 | C15H8O    | 005737-13-3 | 96   |
| 2    |    | 1,2,4,8-Tetramethylbicyclo[6.3.0...] | 204 | C15H24    | 137235-51-9 | 55   |
| 3    |    | 6-Phenyl-2-thiouracil                | 204 | C10H8N2OS | 005590-94-3 | 53   |
| 4    |    | 4,5,6,7-Tetrafluoro-2-methyl-1H-...  | 204 | C8H4F4N2  | 081430-75-3 | 53   |
| 5    |    | [1,1'-Biphenyl]-4,4'-dicarbonitrile  | 204 | C14H8N2   | 001591-30-6 | 53   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

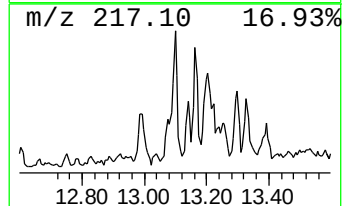
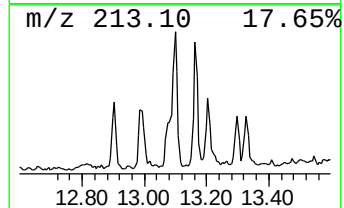
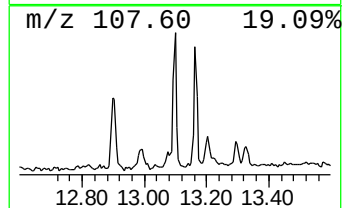
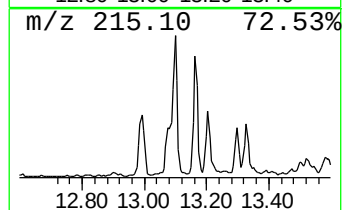
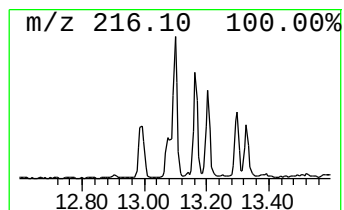
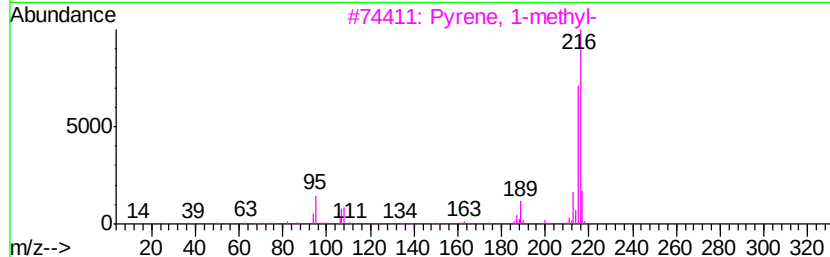
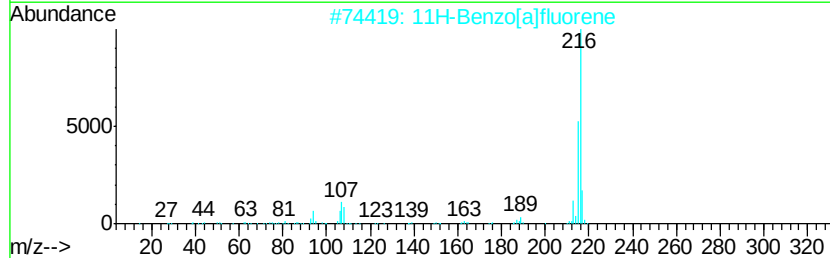
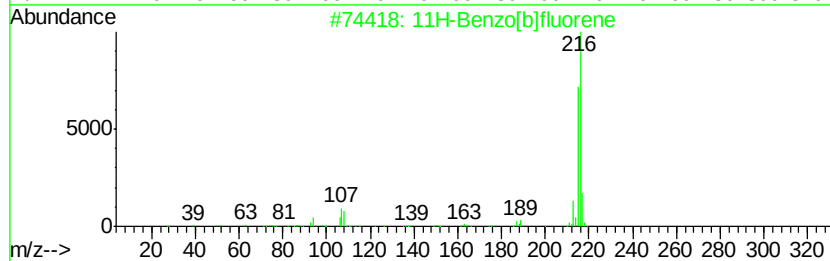
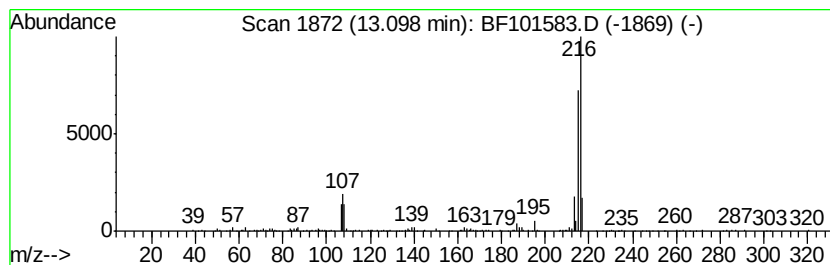
Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.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 11H-Benzo[b]fluorene Concentration Rank 10

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.    |             |      |
|---------|---------|-------------------------|------------------|---------|-------------|------|
| 13.10   | 2.91 ng | 240196                  | Chrysene-d12     | 13.97   |             |      |
| Hit# of | 5       | Tentative ID            | MW               | MolForm | CAS#        | Qual |
| 1       |         | 11H-Benzo[b]fluorene    | 216              | C17H12  | 000243-17-4 | 93   |
| 2       |         | 11H-Benzo[a]fluorene    | 216              | C17H12  | 000238-84-6 | 90   |
| 3       |         | Pyrene, 1-methyl-       | 216              | C17H12  | 002381-21-7 | 90   |
| 4       |         | Fluoranthene, 2-methyl- | 216              | C17H12  | 033543-31-6 | 81   |
| 5       |         | Pyrene, 2-methyl-       | 216              | C17H12  | 003442-78-2 | 80   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

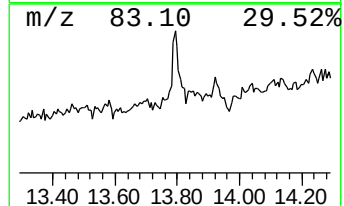
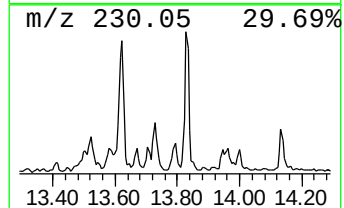
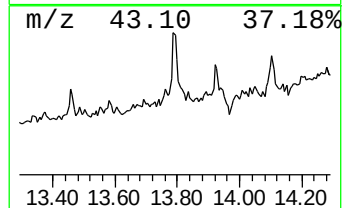
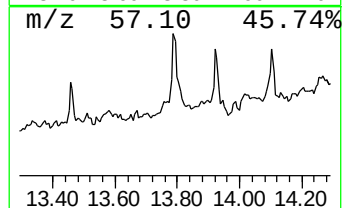
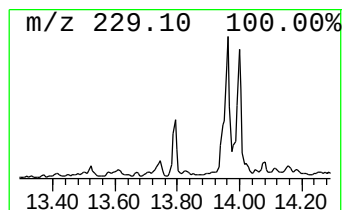
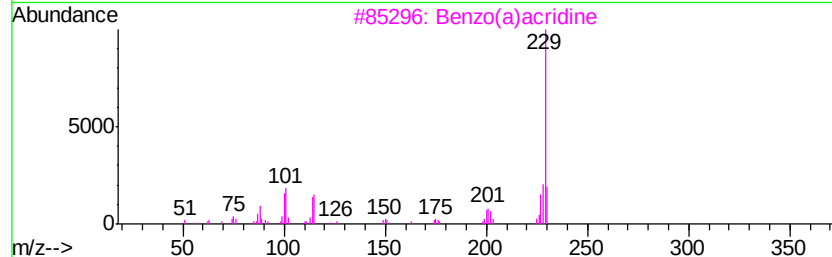
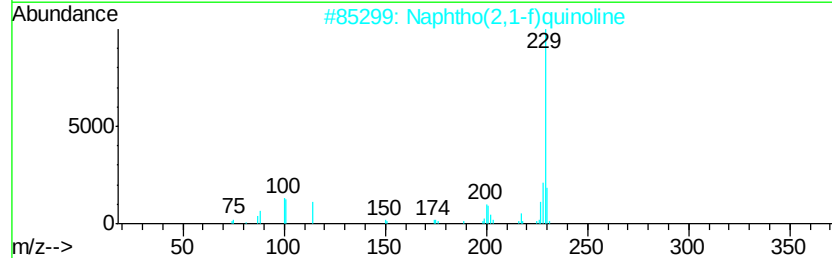
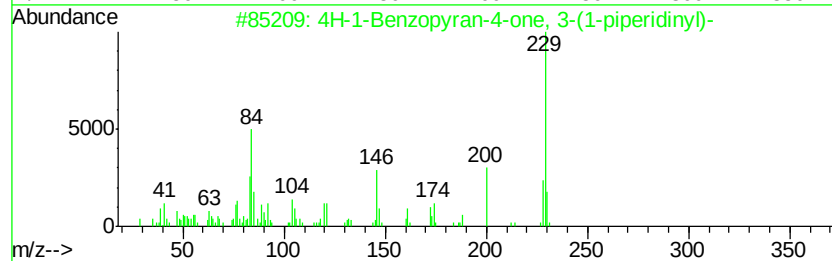
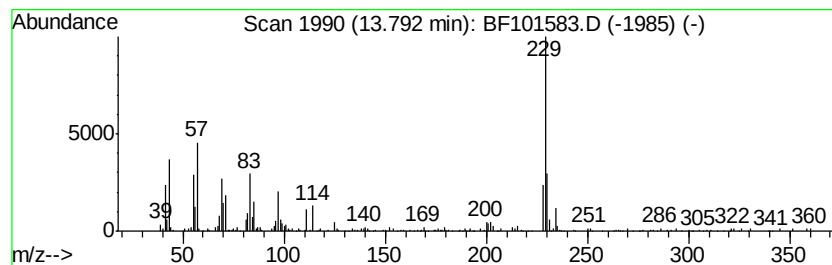
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown13.79 Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.79 | 4.47 ng | 369409 | Chrysene-d12     | 13.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | 4H-1-Benzopyran-4-one, 3-(1-pipe... | 229 | C14H15N02    | 040302-79-2  | 42   |
| 2    |    | Naphtho(2,1-f)quinoline             | 229 | C17H11N      | 000218-08-6  | 38   |
| 3    |    | Benzo(a)acridine                    | 229 | C17H11N      | 000225-11-6  | 37   |
| 4    |    | Benz[c]acridine                     | 229 | C17H11N      | 000225-51-4  | 25   |
| 5    |    | 1-Dichloromethyldimethylsilyloxy... | 312 | C14H30Cl2OSi | 1000283-10-7 | 25   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
Data File : BF101583.D  
Acq On : 20 Dec 2017 22:14  
Operator : SJ/JU  
Sample : I6981-04 2X  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

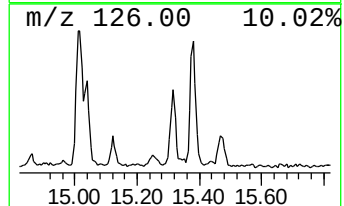
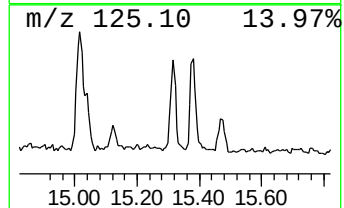
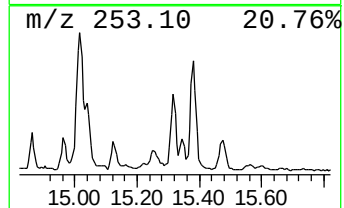
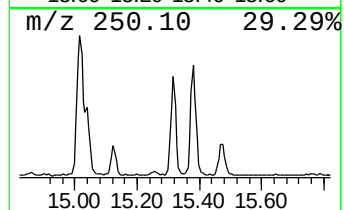
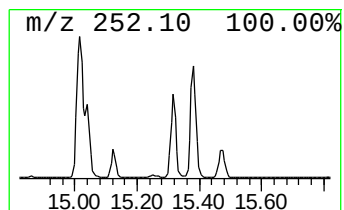
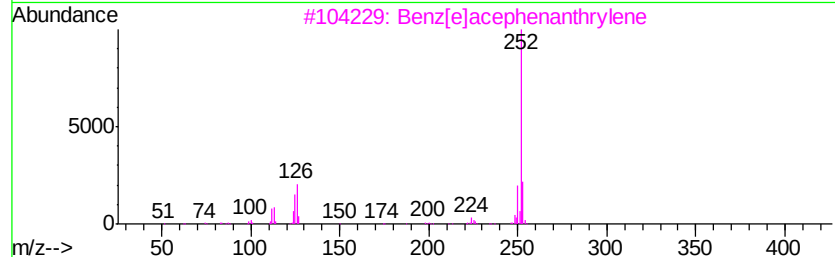
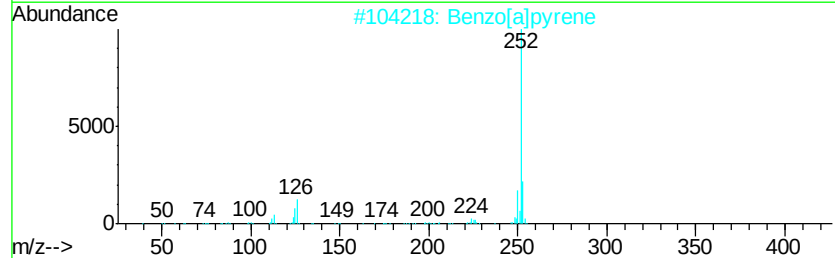
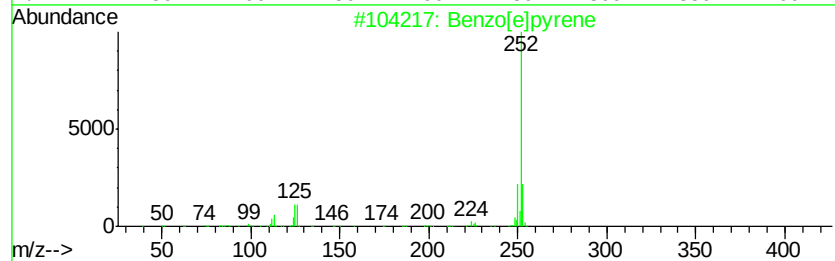
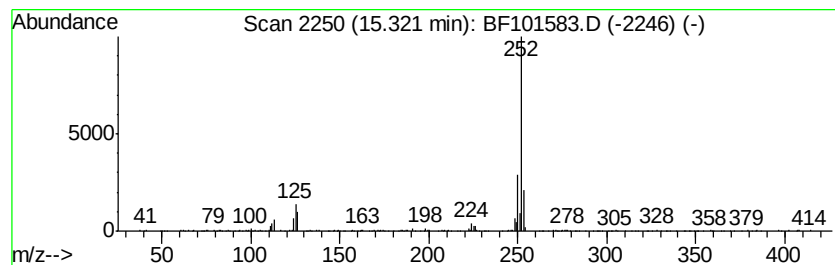
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Benzo[e]pyrene Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.32 | 14.16 ng | 307443 | Perylene-d12     | 15.44 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[e]pyrene           | 252 | C20H12  | 000192-97-2 | 98   |
| 2    |    | Benzo[a]pyrene           | 252 | C20H12  | 000050-32-8 | 96   |
| 3    |    | Benz[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 94   |
| 4    |    | Perylene                 | 252 | C20H12  | 000198-55-0 | 90   |
| 5    |    | Benzo[k]fluoranthene     | 252 | C20H12  | 000207-08-9 | 89   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF122017\  
 Data File : BF101583.D  
 Acq On : 20 Dec 2017 22:14  
 Operator : SJ/JU  
 Sample : I6981-04 2X  
 Misc :  
 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 PHMHA-WCD

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF121417.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pentanone, 4-hy... | 5.00  | 5.0     | ng    | 232910   | 1                     | 6.79  | 938757  | 20.0 |
| unknown6.56          | 6.56  | 39.0    | ng    | 1829950  | 1                     | 6.79  | 938757  | 20.0 |
| Dodecane             | 10.25 | 2.7     | ng    | 152356   | 3                     | 9.83  | 1130660 | 20.0 |
| Hexadecane           | 10.72 | 4.1     | ng    | 188034   | 4                     | 11.32 | 908091  | 20.0 |
| 9H-Fluorene, 2-me... | 10.93 | 2.3     | ng    | 103887   | 4                     | 11.32 | 908091  | 20.0 |
| Octacosane           | 11.17 | 2.6     | ng    | 119558   | 4                     | 11.32 | 908091  | 20.0 |
| Phenanthrene, 2-m... | 11.82 | 3.6     | ng    | 163848   | 4                     | 11.32 | 908091  | 20.0 |
| 4H-Cyclopenta[def... | 11.94 | 5.7     | ng    | 260160   | 4                     | 11.32 | 908091  | 20.0 |
| 1H-Cyclopropa[l]p... | 11.96 | 2.3     | ng    | 104997   | 4                     | 11.32 | 908091  | 20.0 |
| Naphthalene, 2-ph... | 12.12 | 2.5     | ng    | 113773   | 4                     | 11.32 | 908091  | 20.0 |
| Anthracene, 2-ethyl- | 12.27 | 2.3     | ng    | 102067   | 4                     | 11.32 | 908091  | 20.0 |
| Phenanthrene, 1,7... | 12.37 | 2.3     | ng    | 104563   | 4                     | 11.32 | 908091  | 20.0 |
| Cyclopenta(def)ph... | 12.47 | 2.4     | ng    | 108496   | 4                     | 11.32 | 908091  | 20.0 |
| 11H-Benzo[b]fluorene | 13.10 | 2.9     | ng    | 240196   | 5                     | 13.97 | 1651980 | 20.0 |
| unknown13.79         | 13.79 | 4.5     | ng    | 369409   | 5                     | 13.97 | 1651980 | 20.0 |
| Benzo[e]pyrene       | 15.32 | 14.2    | ng    | 307443   | 6                     | 15.44 | 434379  | 20.0 |