

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040318\  
 Data File : BF104198.D  
 Acq On : 3 Apr 2018 23:42  
 Operator : JU/SJ  
 Sample : J2182-17  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 6-WM-DIP-7-WW(1-5)

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-BF032818.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         | 2.263       | 26            | 30          | 48           | rVB      | 86741          | 133741        | 2.26%           | 0.194%        |
| 2         | 5.198       | 525           | 529         | 539          | rVB      | 56955          | 80087         | 1.35%           | 0.116%        |
| 3         | 5.587       | 592           | 595         | 621          | rBV      | 1287646        | 1927061       | 32.53%          | 2.792%        |
| 4         | 6.592       | 763           | 766         | 786          | rBV      | 1415455        | 2157824       | 36.42%          | 3.126%        |
| 5         | 6.734       | 786           | 790         | 824          | rVB      | 1729069        | 2381406       | 40.19%          | 3.450%        |
| 6         | 6.957       | 825           | 828         | 844          | rBV      | 382912         | 547929        | 9.25%           | 0.794%        |
| 7         | 7.110       | 851           | 854         | 869          | rBV      | 1493535        | 1565880       | 26.43%          | 2.268%        |
| 8         | 7.522       | 921           | 924         | 947          | rBV      | 813897         | 1343646       | 22.68%          | 1.947%        |
| 9         | 8.239       | 1042          | 1046        | 1048         | rBV      | 750952         | 714561        | 12.06%          | 1.035%        |
| 10        | 8.263       | 1048          | 1050        | 1072         | rVB      | 519496         | 701826        | 11.85%          | 1.017%        |
| 11        | 8.951       | 1164          | 1167        | 1180         | rBV      | 325373         | 405290        | 6.84%           | 0.587%        |
| 12        | 9.051       | 1180          | 1184        | 1196         | rVB      | 269383         | 290097        | 4.90%           | 0.420%        |
| 13        | 9.310       | 1224          | 1228        | 1237         | rBV      | 2349633        | 2378447       | 40.15%          | 3.446%        |
| 14        | 9.375       | 1237          | 1239        | 1243         | rVV      | 84831          | 105784        | 1.79%           | 0.153%        |
| 15        | 9.416       | 1243          | 1246        | 1256         | rVV      | 89755          | 154968        | 2.62%           | 0.225%        |
| 16        | 9.510       | 1260          | 1262        | 1269         | rVV      | 45473          | 69651         | 1.18%           | 0.101%        |
| 17        | 9.581       | 1269          | 1274        | 1282         | rVV      | 110787         | 193910        | 3.27%           | 0.281%        |
| 18        | 9.651       | 1282          | 1286        | 1289         | rVV      | 192508         | 216668        | 3.66%           | 0.314%        |
| 19        | 9.681       | 1289          | 1291        | 1292         | rVV      | 128298         | 111815        | 1.89%           | 0.162%        |
| 20        | 9.704       | 1292          | 1295        | 1301         | rVV      | 213520         | 265392        | 4.48%           | 0.384%        |
| 21        | 9.769       | 1301          | 1306        | 1314         | rVV      | 88470          | 141230        | 2.38%           | 0.205%        |
| 22        | 9.857       | 1318          | 1321        | 1326         | rVV2     | 78882          | 95440         | 1.61%           | 0.138%        |
| 23        | 9.904       | 1326          | 1329        | 1333         | rVB      | 112994         | 89102         | 1.50%           | 0.129%        |
| 24        | 9.998       | 1342          | 1345        | 1348         | rBV      | 971161         | 864158        | 14.59%          | 1.252%        |
| 25        | 10.028      | 1348          | 1350        | 1357         | rVB      | 893757         | 836016        | 14.11%          | 1.211%        |
| 26        | 10.151      | 1366          | 1371        | 1375         | rVB3     | 72784          | 96861         | 1.63%           | 0.140%        |
| 27        | 10.204      | 1375          | 1380        | 1384         | rBV      | 407146         | 527169        | 8.90%           | 0.764%        |
| 28        | 10.310      | 1393          | 1398        | 1401         | rBV2     | 51663          | 73970         | 1.25%           | 0.107%        |
| 29        | 10.404      | 1408          | 1414        | 1421         | rBV      | 167020         | 184836        | 3.12%           | 0.268%        |
| 30        | 10.551      | 1434          | 1439        | 1445         | rBV      | 1194550        | 1267218       | 21.39%          | 1.836%        |
| 31        | 10.628      | 1448          | 1452        | 1456         | rVV4     | 163995         | 315401        | 5.32%           | 0.457%        |
| 32        | 10.680      | 1459          | 1461        | 1463         | rVV2     | 74393          | 84790         | 1.43%           | 0.123%        |
| 33        | 10.704      | 1463          | 1465        | 1470         | rVB      | 157251         | 154445        | 2.61%           | 0.224%        |
| 34        | 10.792      | 1476          | 1480        | 1486         | rBV2     | 1361015        | 1633316       | 27.57%          | 2.366%        |

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 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 6-WM-DIP-7-WW(1-5)

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-BF032818.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 35 | 10.880 | 1492 | 1495 | 1505 | rVB3 | 135672  | 230049  | 3.88%   | 0.333% |
| 36 | 10.980 | 1509 | 1512 | 1517 | rBV  | 105127  | 120748  | 2.04%   | 0.175% |
| 37 | 11.075 | 1524 | 1528 | 1529 | rBV  | 88164   | 97070   | 1.64%   | 0.141% |
| 38 | 11.098 | 1529 | 1532 | 1535 | rVV  | 179095  | 226049  | 3.82%   | 0.327% |
| 39 | 11.127 | 1535 | 1537 | 1541 | rVV2 | 109288  | 128878  | 2.18%   | 0.187% |
| 40 | 11.180 | 1543 | 1546 | 1548 | rVV  | 92232   | 82692   | 1.40%   | 0.120% |
| 41 | 11.222 | 1548 | 1553 | 1555 | rVV4 | 55118   | 84308   | 1.42%   | 0.122% |
| 42 | 11.245 | 1555 | 1557 | 1563 | rVB2 | 73079   | 80760   | 1.36%   | 0.117% |
| 43 | 11.351 | 1569 | 1575 | 1579 | rBV3 | 550594  | 625745  | 10.56%  | 0.907% |
| 44 | 11.392 | 1579 | 1582 | 1589 | rVB  | 314865  | 364806  | 6.16%   | 0.528% |
| 45 | 11.498 | 1596 | 1600 | 1601 | rBV  | 681379  | 690861  | 11.66%  | 1.001% |
| 46 | 11.533 | 1601 | 1606 | 1611 | rVV  | 4098968 | 5924635 | 100.00% | 8.583% |
| 47 | 11.575 | 1611 | 1613 | 1634 | rVV  | 1425106 | 2131215 | 35.97%  | 3.087% |
| 48 | 11.733 | 1634 | 1640 | 1646 | rVV  | 509551  | 804044  | 13.57%  | 1.165% |
| 49 | 11.780 | 1646 | 1648 | 1651 | rVV  | 120788  | 143077  | 2.41%   | 0.207% |
| 50 | 11.827 | 1654 | 1656 | 1667 | rVV4 | 73350   | 166672  | 2.81%   | 0.241% |
| 51 | 11.904 | 1667 | 1669 | 1675 | rVB2 | 47386   | 74342   | 1.25%   | 0.108% |
| 52 | 11.998 | 1681 | 1685 | 1687 | rBV  | 558622  | 575151  | 9.71%   | 0.833% |
| 53 | 12.027 | 1687 | 1690 | 1696 | rVV  | 758573  | 804499  | 13.58%  | 1.165% |
| 54 | 12.074 | 1696 | 1698 | 1701 | rVV  | 226648  | 219949  | 3.71%   | 0.319% |
| 55 | 12.116 | 1701 | 1705 | 1712 | rVB2 | 1132687 | 1589057 | 26.82%  | 2.302% |
| 56 | 12.292 | 1732 | 1735 | 1739 | rBV  | 376051  | 397023  | 6.70%   | 0.575% |
| 57 | 12.439 | 1758 | 1760 | 1763 | rBV2 | 92801   | 85725   | 1.45%   | 0.124% |
| 58 | 12.474 | 1763 | 1766 | 1768 | rBV  | 80199   | 66754   | 1.13%   | 0.097% |
| 59 | 12.498 | 1768 | 1770 | 1775 | rVB  | 72770   | 83623   | 1.41%   | 0.121% |
| 60 | 12.545 | 1775 | 1778 | 1782 | rBV  | 201796  | 261476  | 4.41%   | 0.379% |
| 61 | 12.586 | 1783 | 1785 | 1791 | rVB2 | 95207   | 141088  | 2.38%   | 0.204% |
| 62 | 12.727 | 1803 | 1809 | 1816 | rBV  | 3562728 | 5255627 | 88.71%  | 7.614% |
| 63 | 12.869 | 1830 | 1833 | 1836 | rBV2 | 138199  | 145143  | 2.45%   | 0.210% |
| 64 | 12.963 | 1841 | 1849 | 1853 | rBV2 | 3179460 | 5230590 | 88.29%  | 7.578% |
| 65 | 12.992 | 1853 | 1854 | 1860 | rVB  | 235045  | 221892  | 3.75%   | 0.321% |
| 66 | 13.080 | 1864 | 1869 | 1873 | rBV2 | 1996362 | 2201141 | 37.15%  | 3.189% |
| 67 | 13.163 | 1879 | 1883 | 1888 | rBV3 | 229760  | 397157  | 6.70%   | 0.575% |
| 68 | 13.280 | 1895 | 1903 | 1910 | rBV  | 817040  | 1261449 | 21.29%  | 1.827% |
| 69 | 13.345 | 1910 | 1914 | 1917 | rVV  | 834592  | 863941  | 14.58%  | 1.252% |
| 70 | 13.380 | 1918 | 1920 | 1922 | rVV2 | 299616  | 338304  | 5.71%   | 0.490% |
| 71 | 13.404 | 1922 | 1924 | 1927 | rVB2 | 210626  | 204835  | 3.46%   | 0.297% |

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

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

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032818.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 72 | 13.474 | 1933 | 1936 | 1938 | rBV  | 181158  | 170116  | 2.87%  | 0.246% |
| 73 | 13.504 | 1938 | 1941 | 1950 | rVB2 | 233893  | 327536  | 5.53%  | 0.475% |
| 74 | 13.627 | 1960 | 1962 | 1964 | rBV2 | 74484   | 65800   | 1.11%  | 0.095% |
| 75 | 13.698 | 1972 | 1974 | 1981 | rVB  | 149036  | 187966  | 3.17%  | 0.272% |
| 76 | 13.757 | 1982 | 1984 | 1987 | rBV  | 89051   | 104444  | 1.76%  | 0.151% |
| 77 | 13.904 | 2006 | 2009 | 2010 | rBV  | 209001  | 211392  | 3.57%  | 0.306% |
| 78 | 13.927 | 2010 | 2013 | 2015 | rVV  | 320115  | 341210  | 5.76%  | 0.494% |
| 79 | 13.951 | 2015 | 2017 | 2024 | rVB4 | 449832  | 656860  | 11.09% | 0.952% |
| 80 | 14.151 | 2046 | 2051 | 2055 | rBV2 | 1583224 | 2868834 | 48.42% | 4.156% |
| 81 | 14.186 | 2055 | 2057 | 2064 | rVV  | 1666709 | 1629245 | 27.50% | 2.360% |
| 82 | 14.263 | 2067 | 2070 | 2076 | rVB  | 367303  | 359535  | 6.07%  | 0.521% |
| 83 | 14.545 | 2116 | 2118 | 2121 | rVB  | 184636  | 152419  | 2.57%  | 0.221% |
| 84 | 14.645 | 2133 | 2135 | 2138 | rVB2 | 124364  | 112850  | 1.90%  | 0.163% |
| 85 | 14.674 | 2138 | 2140 | 2142 | rBV  | 109717  | 108273  | 1.83%  | 0.157% |
| 86 | 15.245 | 2232 | 2237 | 2240 | rBV  | 1004936 | 1788025 | 30.18% | 2.590% |
| 87 | 15.357 | 2253 | 2256 | 2260 | rBV  | 169981  | 208889  | 3.53%  | 0.303% |
| 88 | 15.568 | 2288 | 2292 | 2298 | rBV  | 566352  | 793289  | 13.39% | 1.149% |
| 89 | 15.639 | 2299 | 2304 | 2311 | rVV  | 853158  | 1361591 | 22.98% | 1.973% |
| 90 | 15.704 | 2311 | 2315 | 2318 | rVV  | 326630  | 507047  | 8.56%  | 0.735% |
| 91 | 15.733 | 2318 | 2320 | 2330 | rVB  | 276235  | 428200  | 7.23%  | 0.620% |
| 92 | 16.410 | 2431 | 2435 | 2442 | rVB2 | 170163  | 304078  | 5.13%  | 0.441% |
| 93 | 16.874 | 2510 | 2514 | 2520 | rVB3 | 109487  | 189427  | 3.20%  | 0.274% |
| 94 | 17.298 | 2579 | 2586 | 2597 | rBV2 | 323288  | 773526  | 13.06% | 1.121% |
| 95 | 17.486 | 2613 | 2618 | 2624 | rBV3 | 106765  | 239507  | 4.04%  | 0.347% |
| 96 | 17.786 | 2661 | 2669 | 2685 | rBV3 | 216656  | 707006  | 11.93% | 1.024% |

Sum of corrected areas: 69027375



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BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

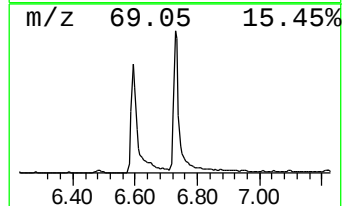
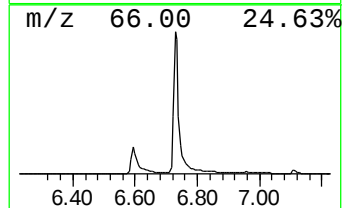
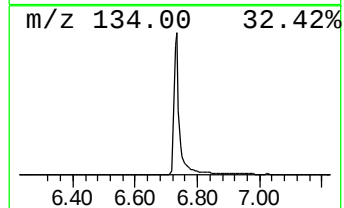
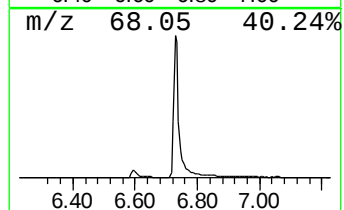
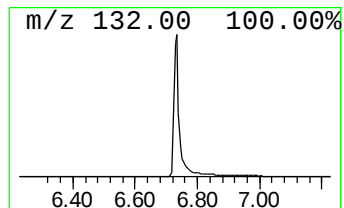
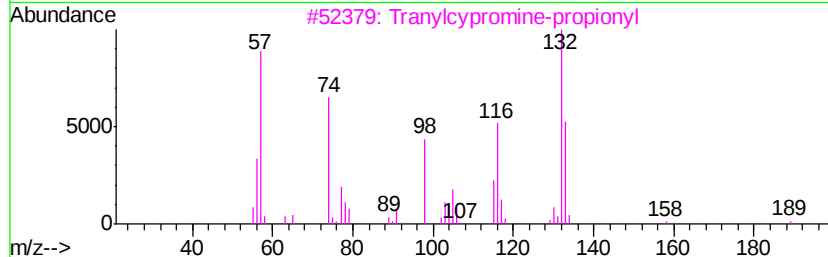
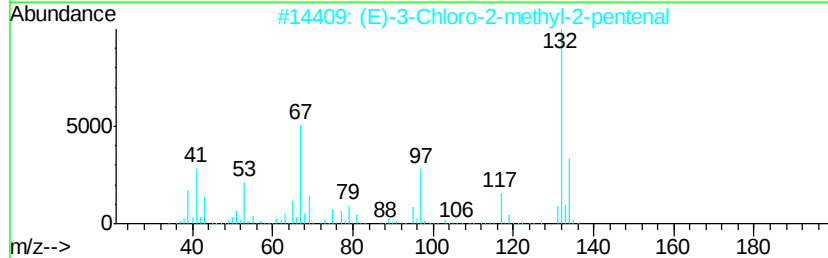
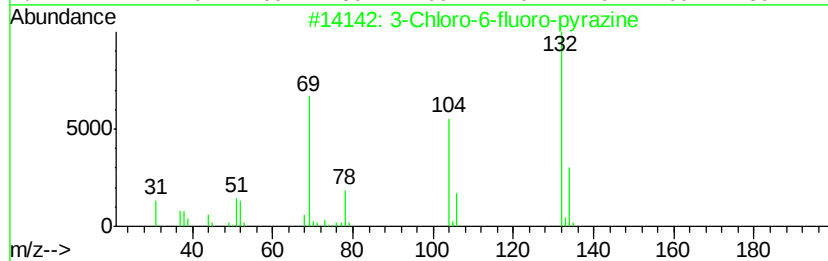
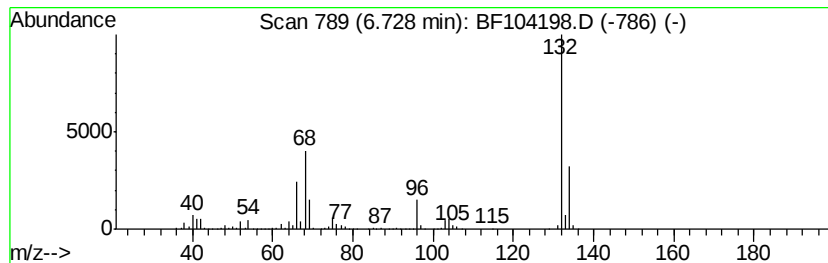
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 unknown6.73 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.73 | 86.92 ng | 2381410 | 1,4-Dichlorobenzene-d4 | 6.96 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 38   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 3    |    | Tranvlcypropine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 4    |    | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2    | 022291-85-6  | 9    |
| 5    |    | 4-Chloro-6-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-01-6  | 9    |



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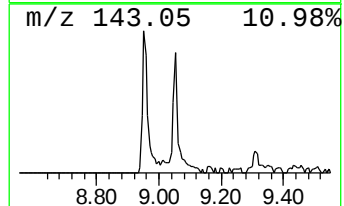
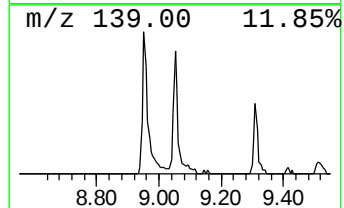
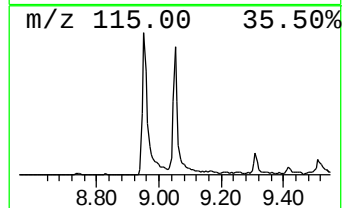
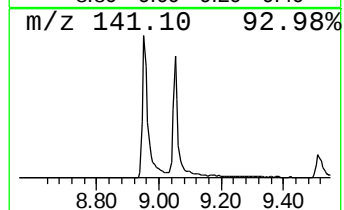
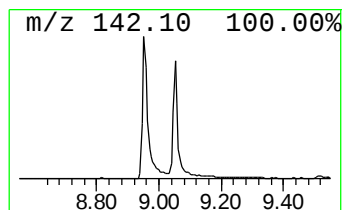
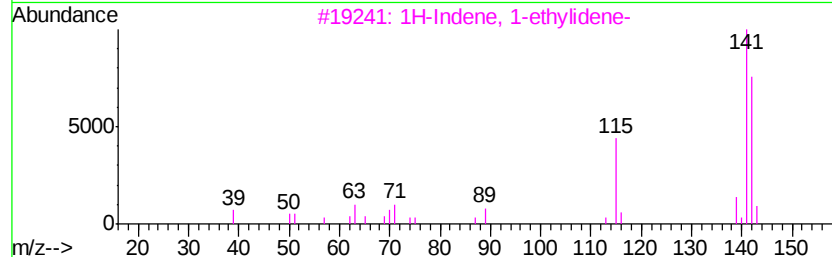
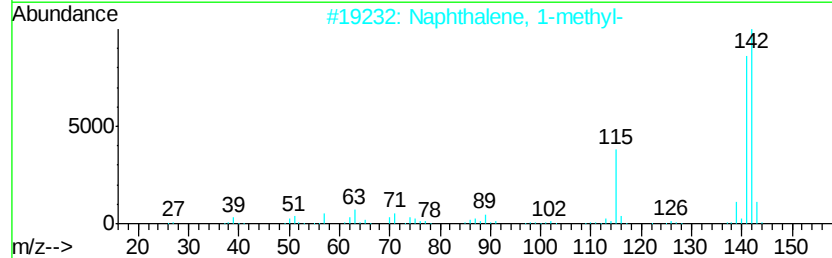
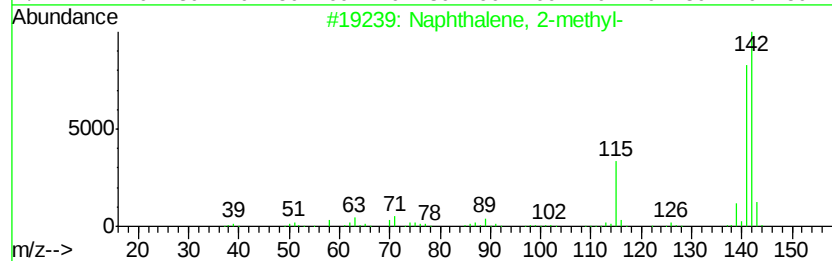
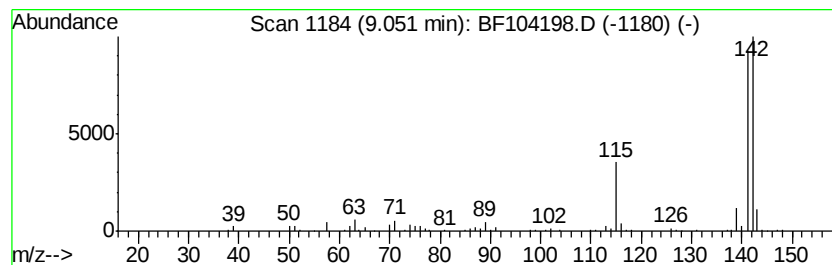
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Naphthalene, 1-methyl- Concentration Rank 14

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.05 | 8.12 ng | 290097 | Naphthalene-d8   | 8.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 97   |
| 2    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 97   |
| 3    |    | 1H-Indene, 1-ethylidene-            | 142 | C11H10  | 002471-83-2 | 94   |
| 4    |    | 1,4-Methanonaphthalene, 1,4-dihy... | 142 | C11H10  | 004453-90-1 | 94   |
| 5    |    | Benzocycloheptatriene               | 142 | C11H10  | 000264-09-5 | 91   |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

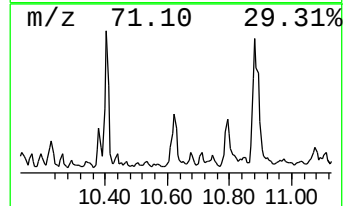
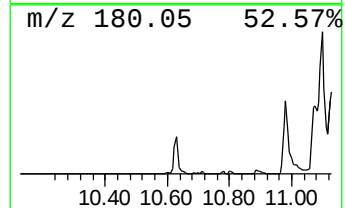
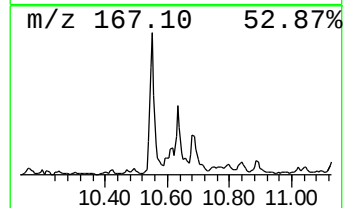
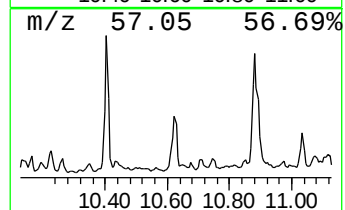
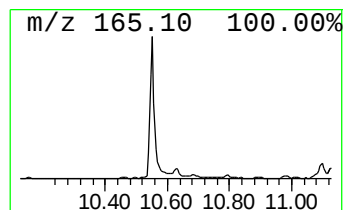
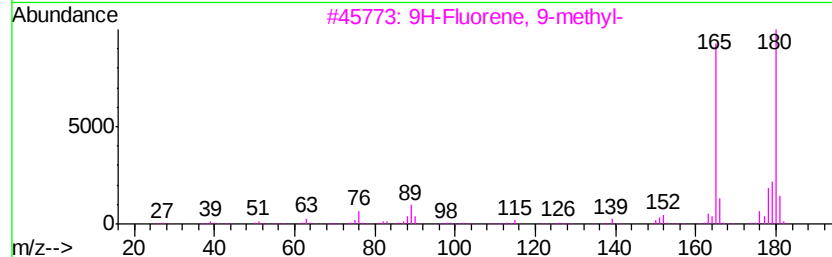
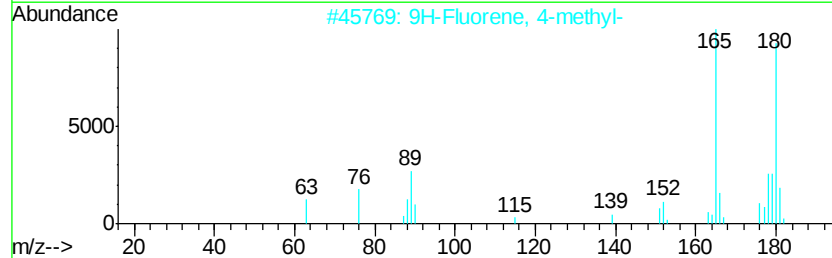
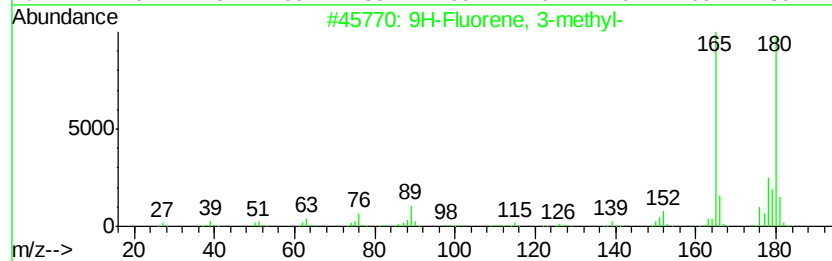
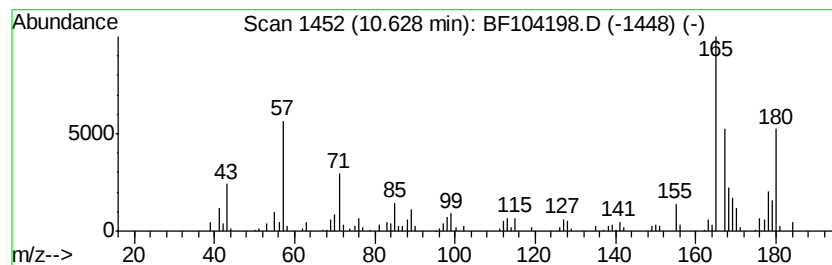
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 unknown10.63 Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.63 | 7.30 ng | 315401 | Acenaphthene-d10 | 10.00 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | 9H-Fluorene, 3-methyl- | 180 | C14H12  | 002523-39-9 | 49   |
| 2    |    | 9H-Fluorene, 4-methyl- | 180 | C14H12  | 001556-99-6 | 45   |
| 3    |    | 9H-Fluorene, 9-methyl- | 180 | C14H12  | 002523-37-7 | 38   |
| 4    |    | 9H-Fluorene, 1-methyl- | 180 | C14H12  | 001730-37-6 | 38   |
| 5    |    | 9H-Fluorene, 2-methyl- | 180 | C14H12  | 001430-97-3 | 38   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

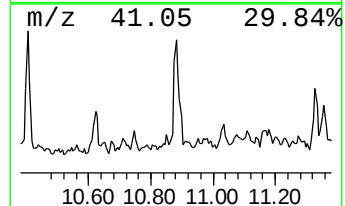
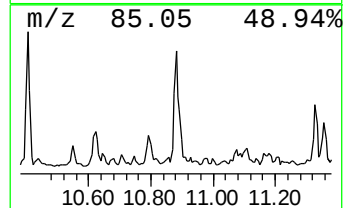
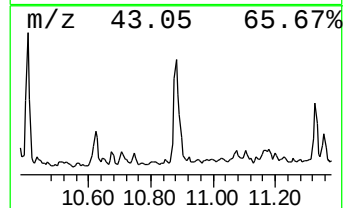
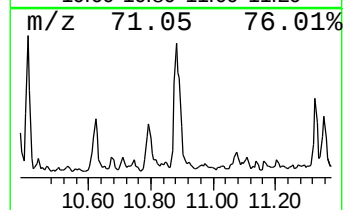
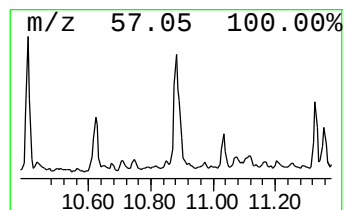
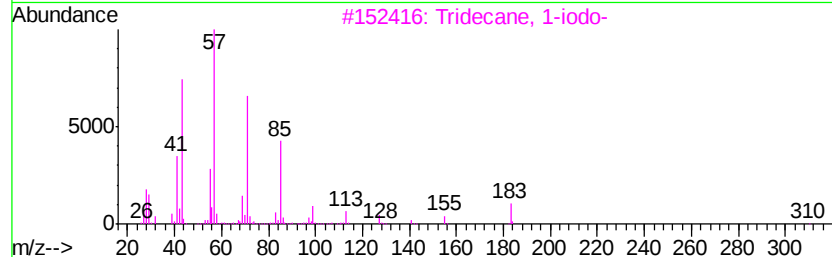
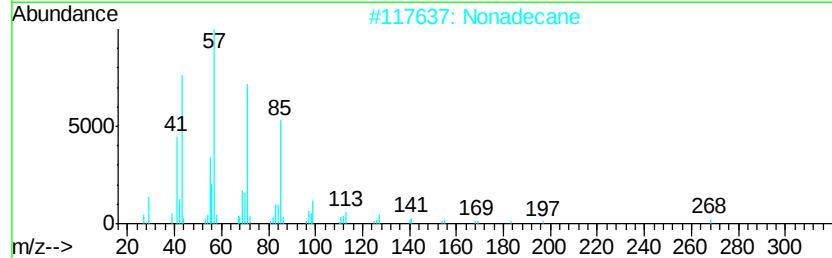
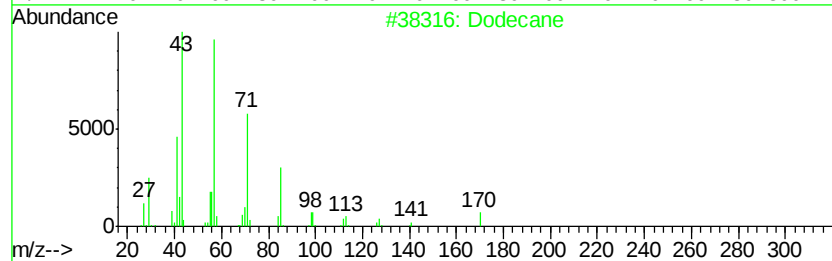
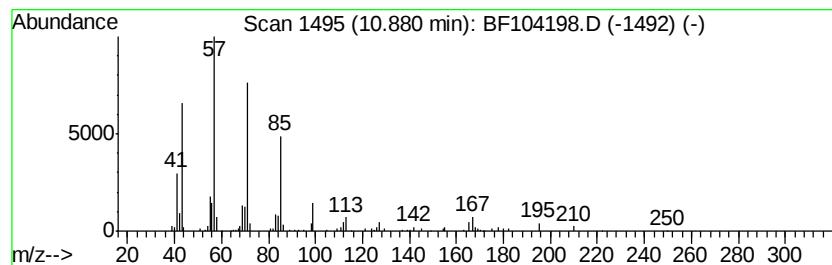
Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Dodecane Concentration Rank 18

| R.T.      | EstConc                | Area   | Relative to ISTD |             | R.T.  |
|-----------|------------------------|--------|------------------|-------------|-------|
| 10.88     | 6.66 ng                | 230049 | Phenanthrene-d10 |             | 11.50 |
| Hit# of 5 | Tentative ID           | MW     | MolForm          | CAS#        | Qual  |
| 1         | Dodecane               | 170    | C12H26           | 000112-40-3 | 93    |
| 2         | Nonadecane             | 268    | C19H40           | 000629-92-5 | 90    |
| 3         | Tridecane, 1-iodo-     | 310    | C13H27I          | 035599-77-0 | 86    |
| 4         | Hexadecane             | 226    | C16H34           | 000544-76-3 | 86    |
| 5         | Heptacosane, 1-chloro- | 414    | C27H55Cl         | 062016-79-9 | 86    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

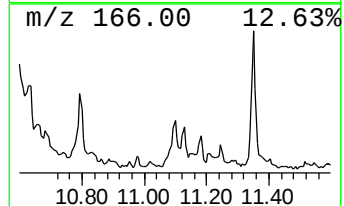
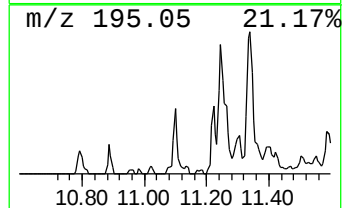
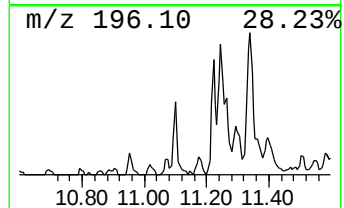
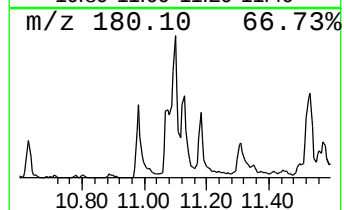
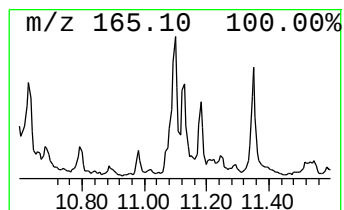
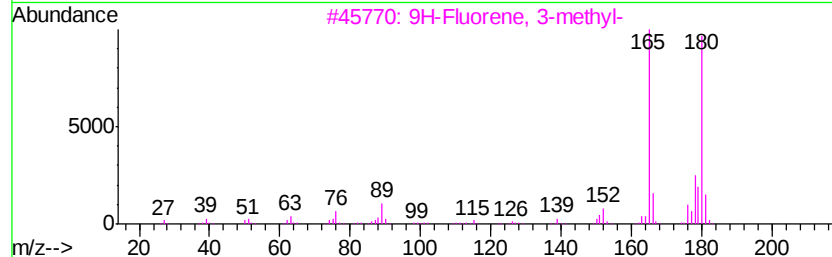
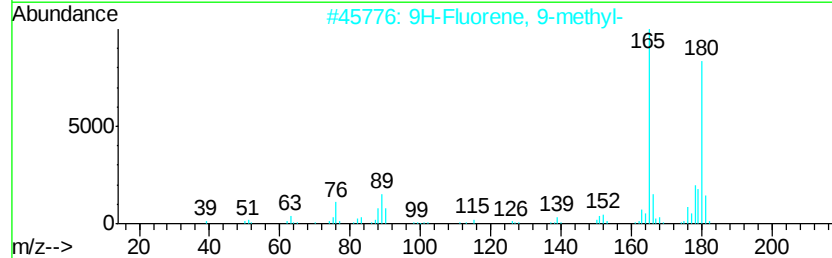
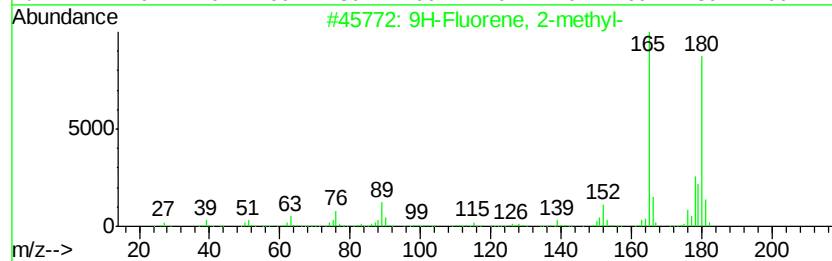
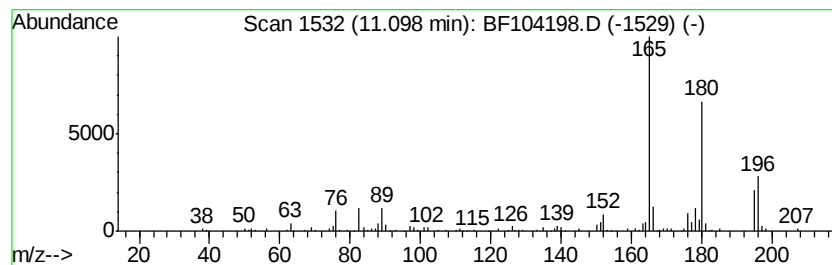
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 9H-Fluorene, 2-methyl- Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.10 | 6.54 ng | 226049 | Phenanthrene-d10 | 11.50 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 9H-Fluorene, 2-methyl-              | 180 | C14H12    | 001430-97-3 | 86   |
| 2    |    | 9H-Fluorene, 9-methyl-              | 180 | C14H12    | 002523-37-7 | 58   |
| 3    |    | 9H-Fluorene, 3-methyl-              | 180 | C14H12    | 002523-39-9 | 58   |
| 4    |    | Benzenamine, N-ethyl-2-methyl-5-... | 180 | C9H12N2O2 | 056288-95-0 | 47   |
| 5    |    | 2',4'-Dimethoxyacetophenone         | 180 | C10H12O3  | 000829-20-9 | 47   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

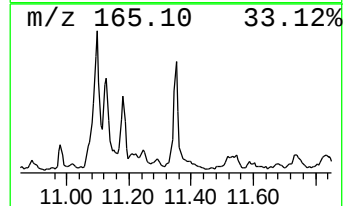
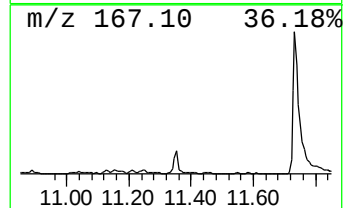
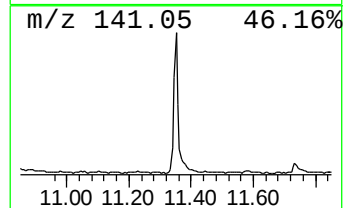
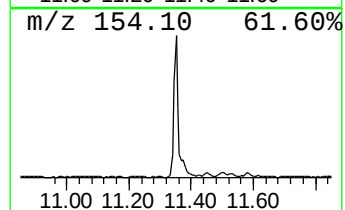
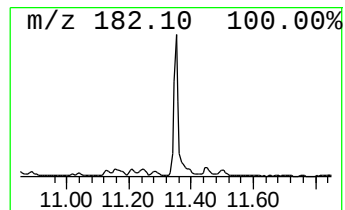
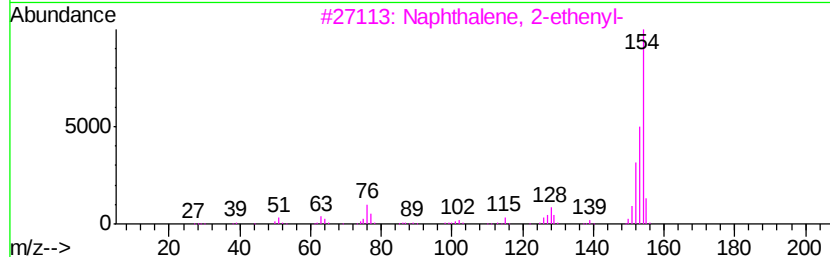
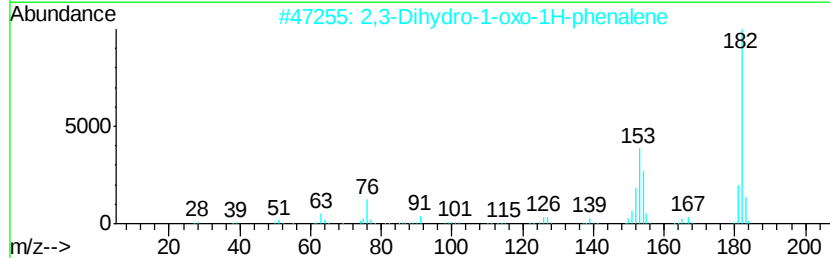
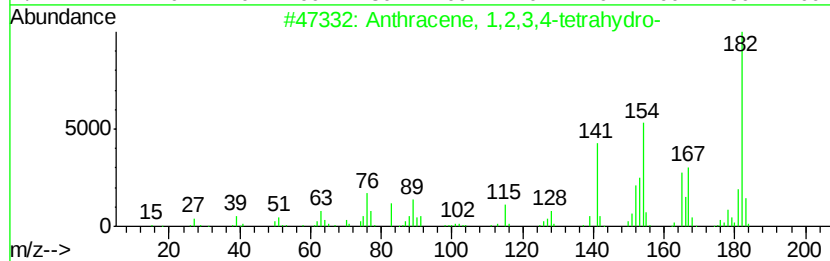
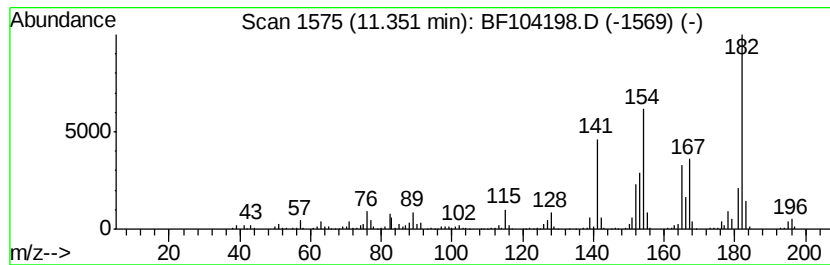
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Anthracene, 1,2,3,4-tetrahy... Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.35 | 18.11 ng | 625745 | Phenanthrene-d10 | 11.50 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 1,2,3,4-tetrahydro-   | 182 | C14H14  | 002141-42-6 | 98   |
| 2    |    |   | 2,3-Dihydro-1-oxo-1H-phenalene    | 182 | C13H10O | 000518-85-4 | 60   |
| 3    |    |   | Naphthalene, 2-ethenyl-           | 154 | C12H10  | 000827-54-3 | 56   |
| 4    |    |   | 3,3'-Dimethylbiphenyl             | 182 | C14H14  | 000612-75-9 | 49   |
| 5    |    |   | Phenanthrene, 1,2,3,4-tetrahydro- | 182 | C14H14  | 001013-08-7 | 49   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

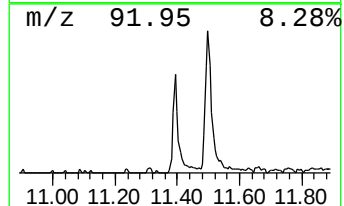
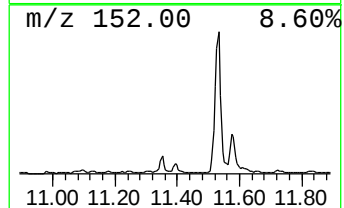
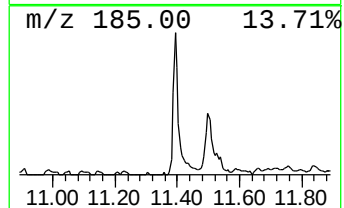
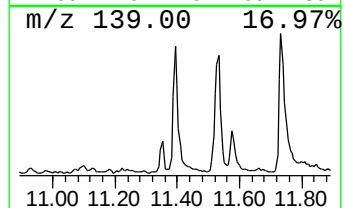
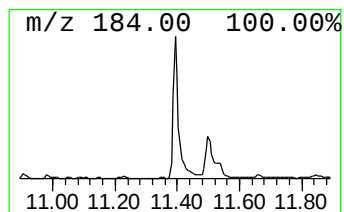
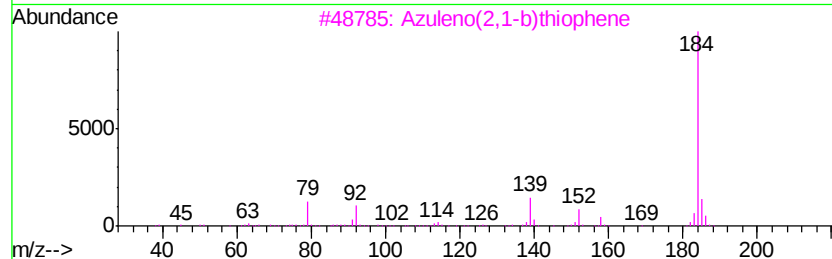
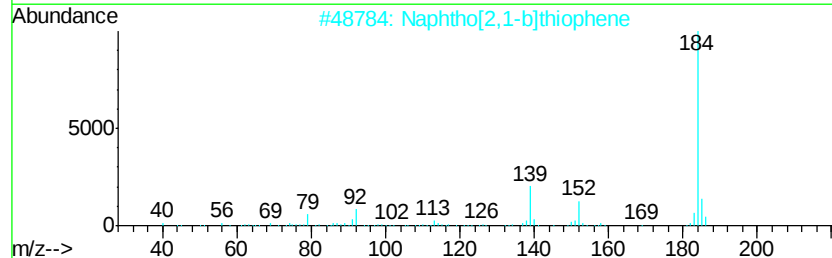
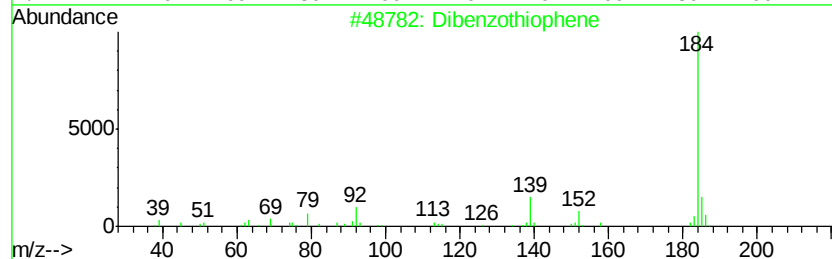
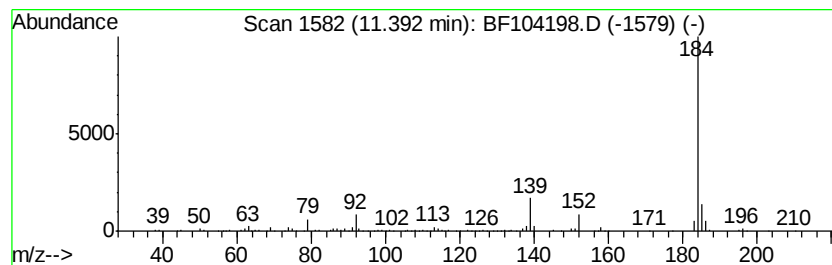
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Dibenzothiophene Concentration Rank 10

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.39 | 10.56 ng | 364806 | Phenanthrene-d10 | 11.50 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Dibenzothiophene                  | 184 | C12H8S  | 000132-65-0 | 96   |
| 2    |    | Naphtho[2,1-b]thiophene           | 184 | C12H8S  | 000233-02-3 | 91   |
| 3    |    | Azuleno(2,1-b)thiophene           | 184 | C12H8S  | 000248-13-5 | 91   |
| 4    |    | Naphtho[2,3-b]thiophene           | 184 | C12H8S  | 000268-77-9 | 87   |
| 5    |    | Naphthalene, 2-ethenyl-6-methoxy- | 184 | C13H12O | 063444-51-9 | 72   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

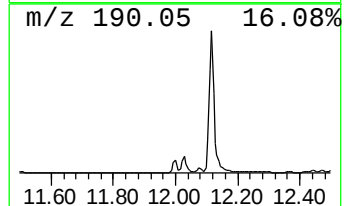
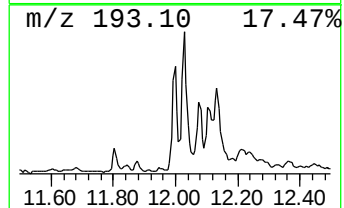
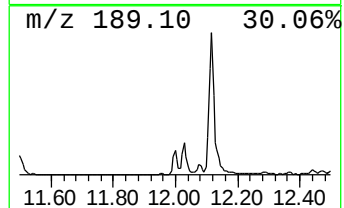
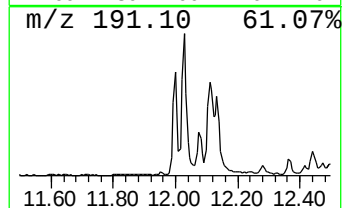
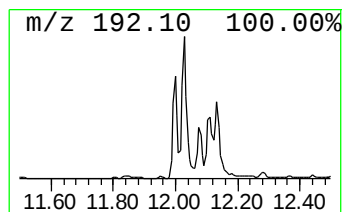
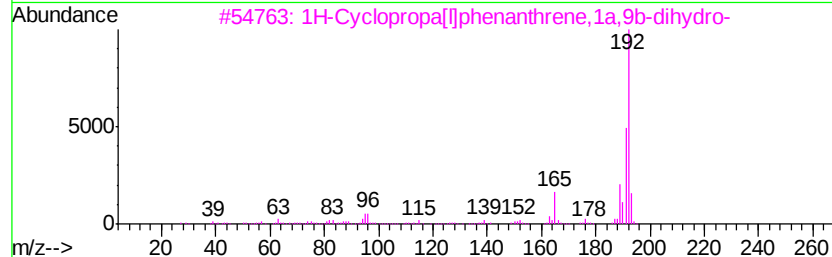
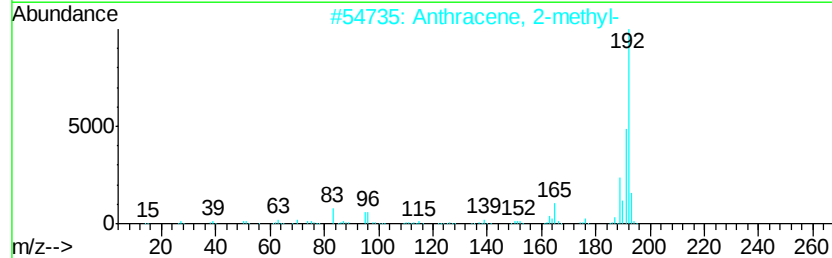
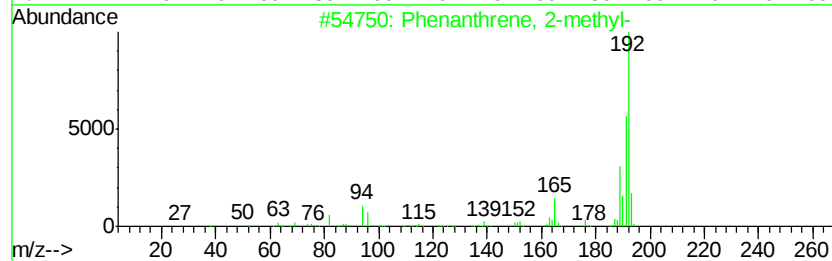
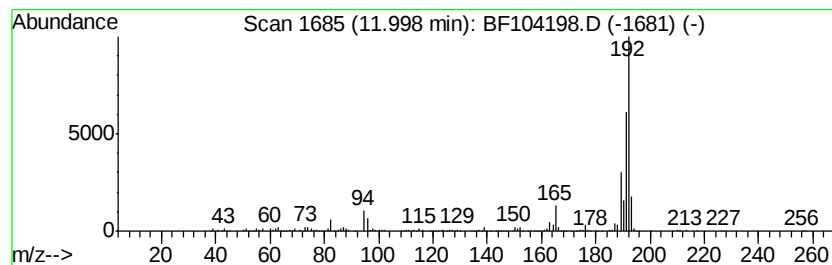
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Phenanthrene, 2-methyl- Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.00 | 16.65 ng | 575151 | Phenanthrene-d10 | 11.50 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 98   |
| 2    |    | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 96   |
| 3    |    | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 95   |
| 4    |    | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 95   |
| 5    |    | Anthracene, 9-methyl-               | 192 | C15H12  | 000779-02-2 | 94   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

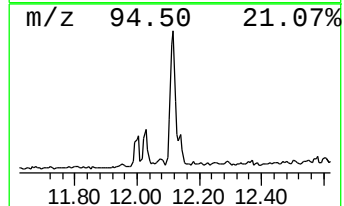
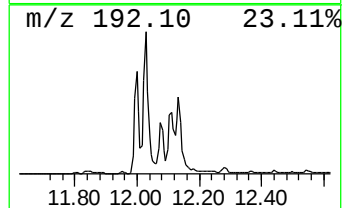
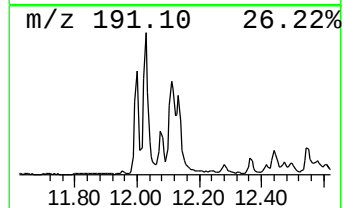
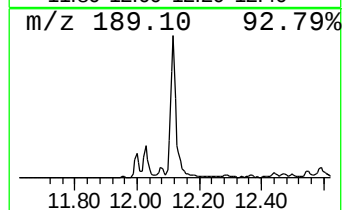
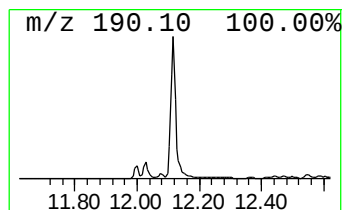
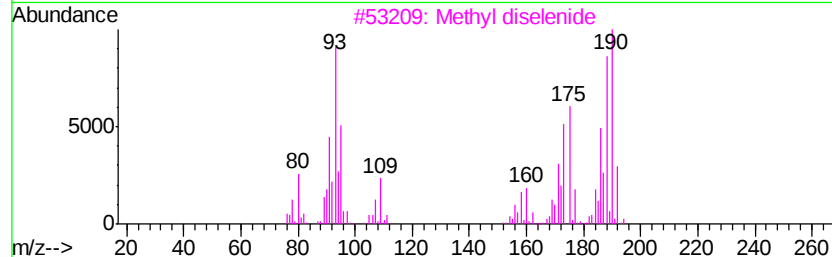
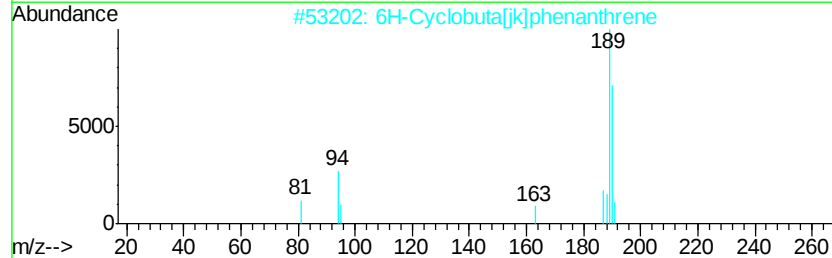
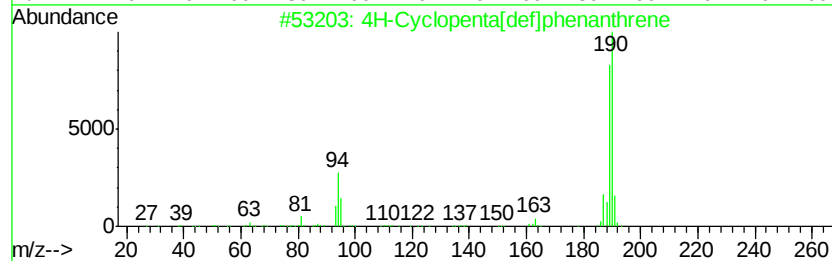
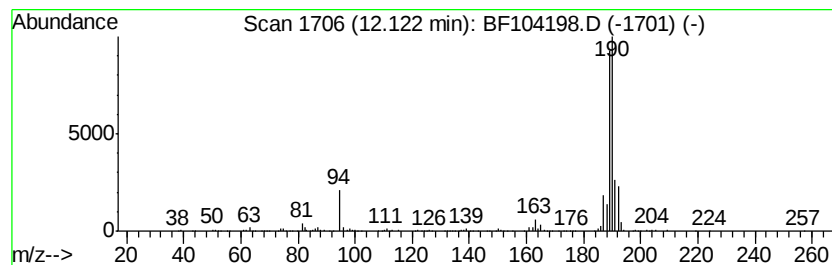
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.12 | 46.00 ng | 1589060 | Phenanthrene-d10 | 11.50 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10   | 000203-64-5 | 94   |
| 2    |    | 6H-Cyclobuta[ik]phenanthrene        | 190 | C15H10   | 083469-43-6 | 64   |
| 3    |    | Methyl diselenide                   | 190 | C2H6Se2  | 007101-31-7 | 45   |
| 4    |    | 2,2'-Bis(4,5-dimethylimidazole)     | 190 | C10H14N4 | 069286-06-2 | 40   |
| 5    |    | 1,4-Naphthalenedione, 2,5-dihydr... | 190 | C10H6O4  | 004923-55-1 | 25   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

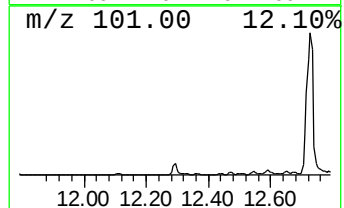
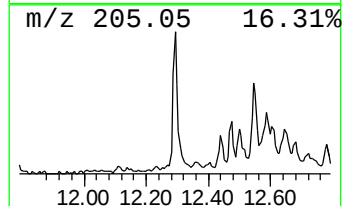
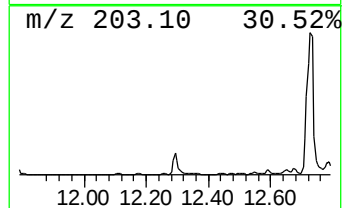
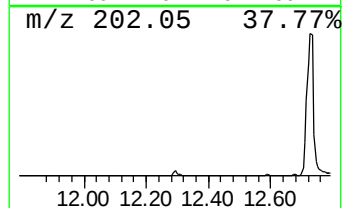
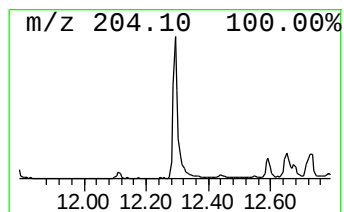
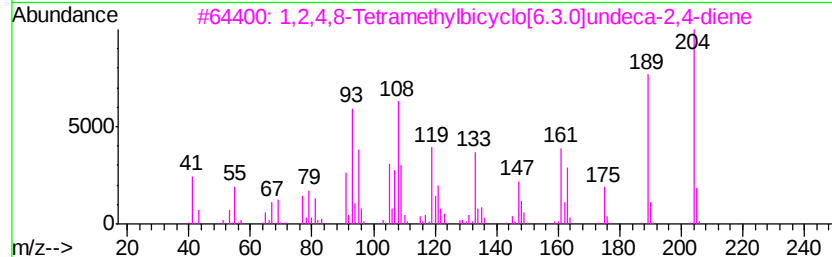
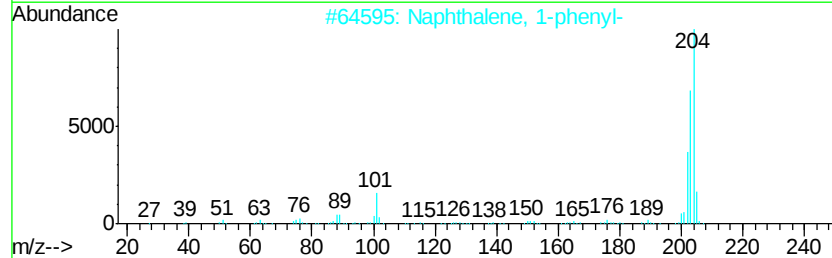
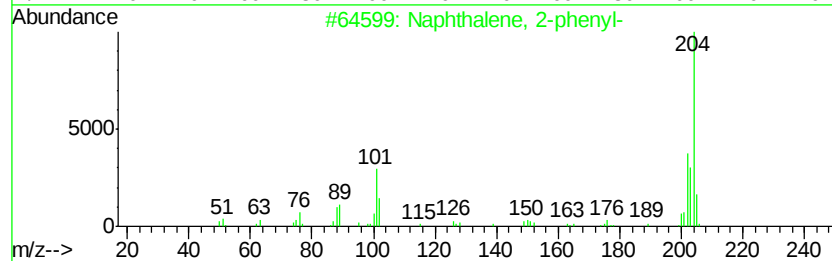
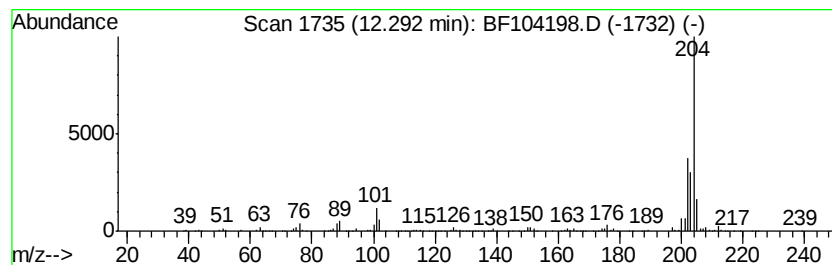
Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 Naphthalene, 2-phenyl- Concentration Rank 9

| R.T.      | EstConc                                                   | Area   | Relative to ISTD | R.T.        |      |
|-----------|-----------------------------------------------------------|--------|------------------|-------------|------|
| 12.29     | 11.49 ng                                                  | 397023 | Phenanthrene-d10 | 11.50       |      |
| Hit# of 5 | Tentative ID                                              | MW     | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 2-phenyl-                                    | 204    | C16H12           | 000612-94-2 | 94   |
| 2         | Naphthalene, 1-phenyl-                                    | 204    | C16H12           | 000605-02-7 | 92   |
| 3         | 1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene         | 204    | C15H24           | 137235-51-9 | 87   |
| 4         | 5,16[1',2']-8,13[1'',2'']-Dibenz[1,2-b:4,5-b']diphenylene | 408    | C32H24           | 005672-97-9 | 87   |
| 5         | Tricyclo[8.2.2.2(4,7)]hexadeca-2,4,6,8-tetraene           | 204    | C16H12           | 006572-60-7 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

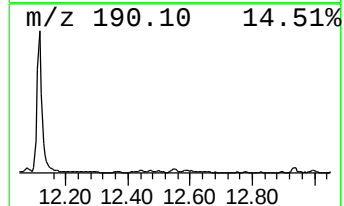
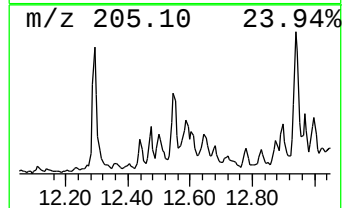
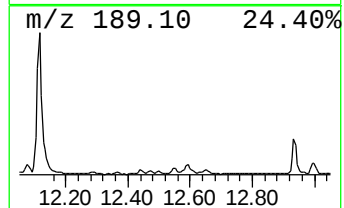
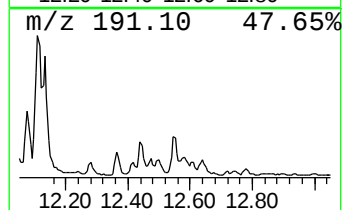
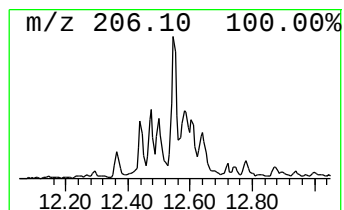
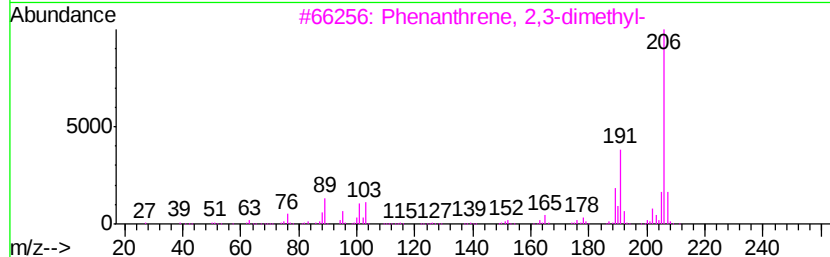
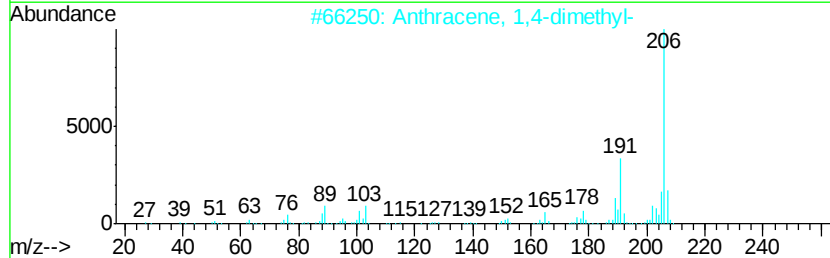
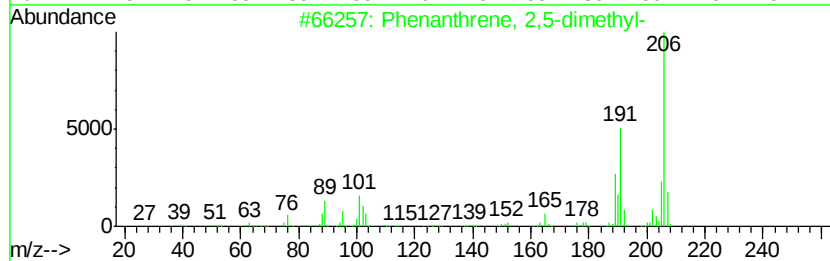
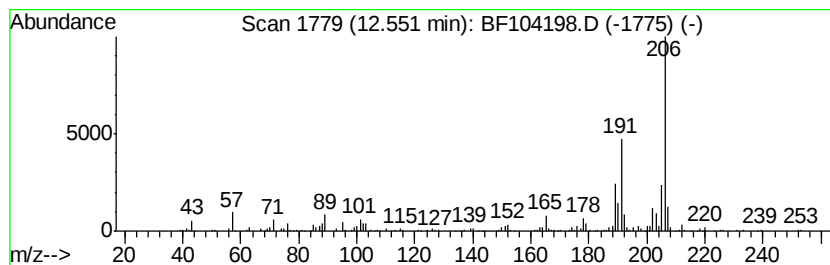
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Phenanthrene, 2,5-dimethyl- Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.55 | 7.57 ng | 261476 | Phenanthrene-d10 | 11.50 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 97   |
| 2    |    |   | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 93   |
| 3    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 91   |
| 4    |    |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 91   |
| 5    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 89   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

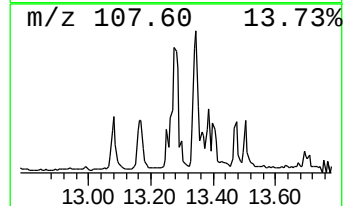
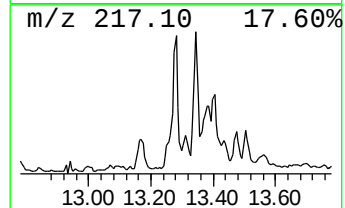
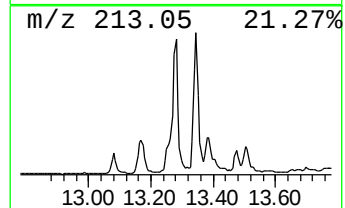
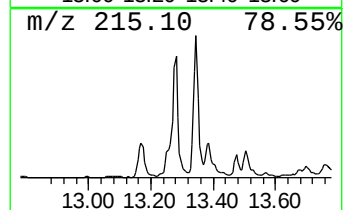
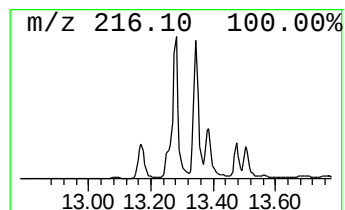
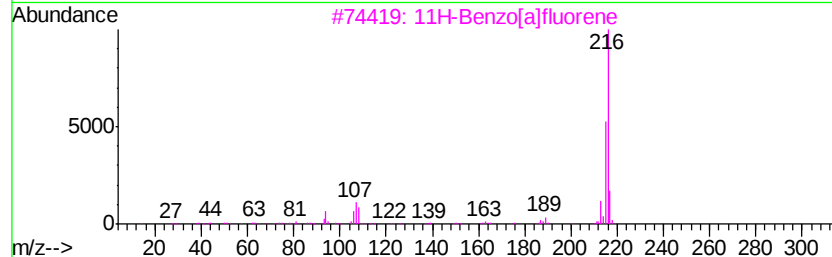
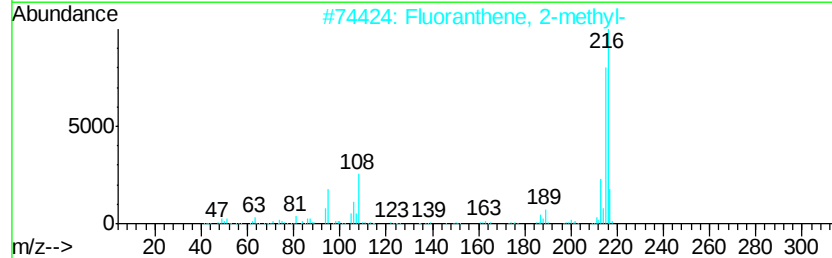
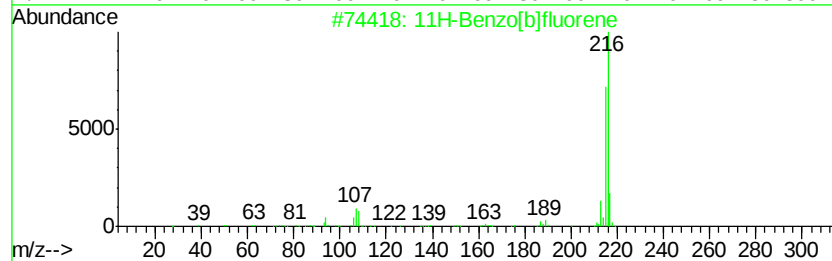
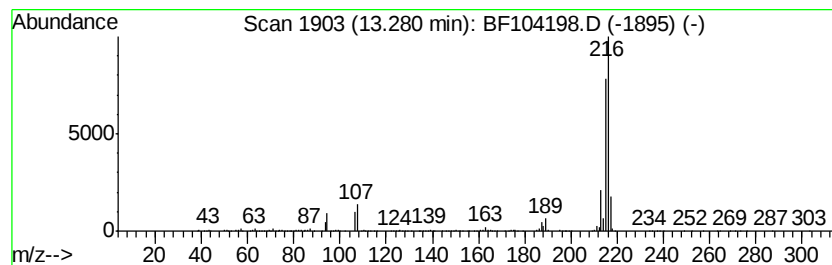
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 12

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 13.28 | 8.79 ng | 1261450 | Chrysene-d12     | 14.16 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 97   |
| 2    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 95   |
| 3    |    | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 94   |
| 4    |    | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 94   |
| 5    |    | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 86   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

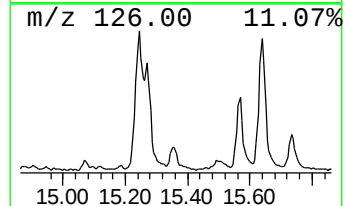
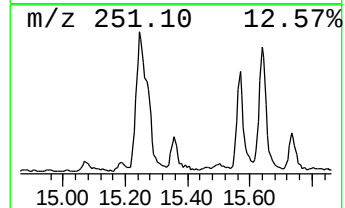
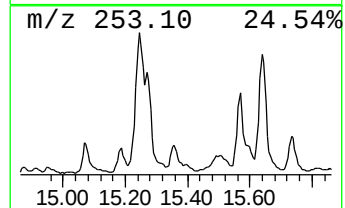
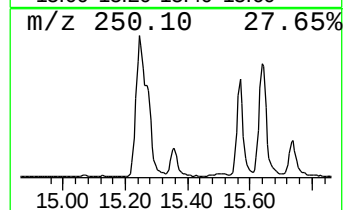
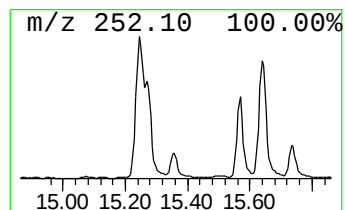
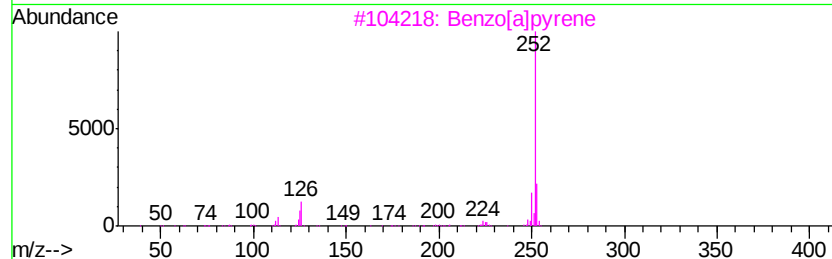
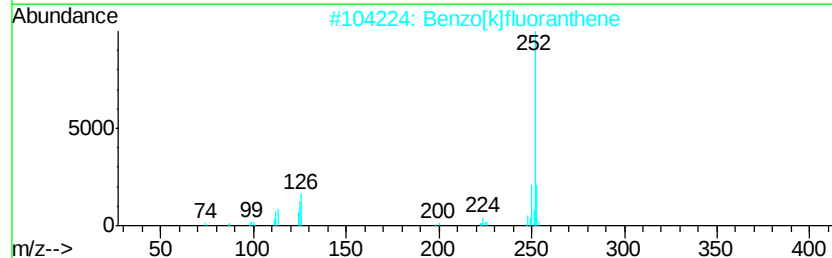
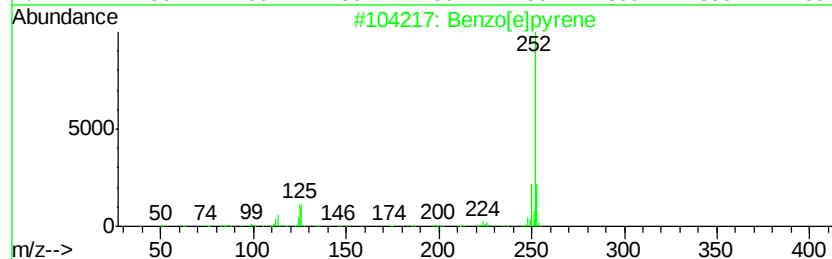
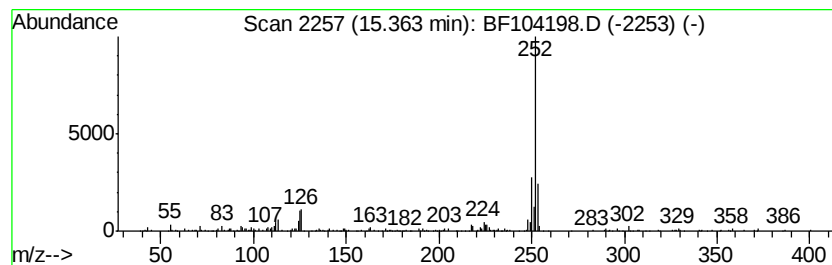
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzo[e]pyrene Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.36 | 8.24 ng | 208889 | Perylene-d12     | 15.70 |

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

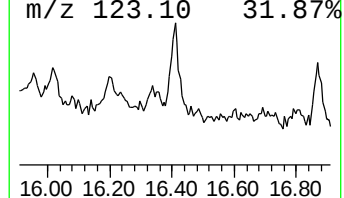
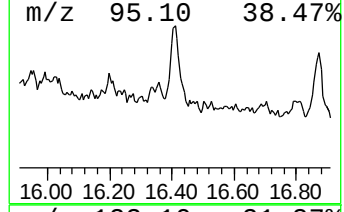
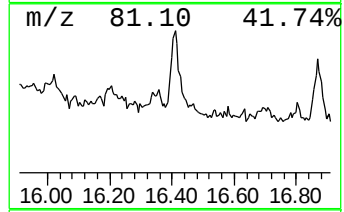
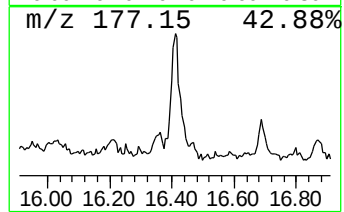
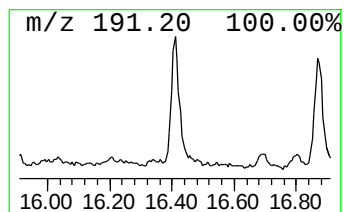
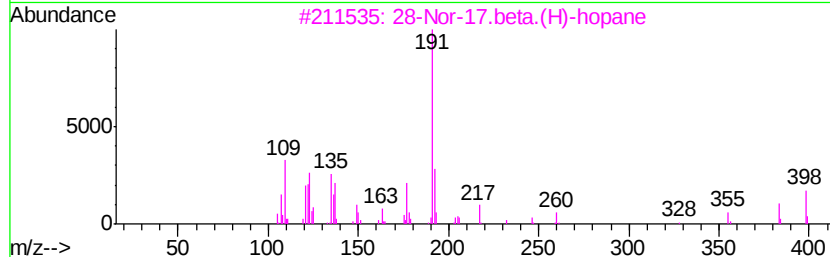
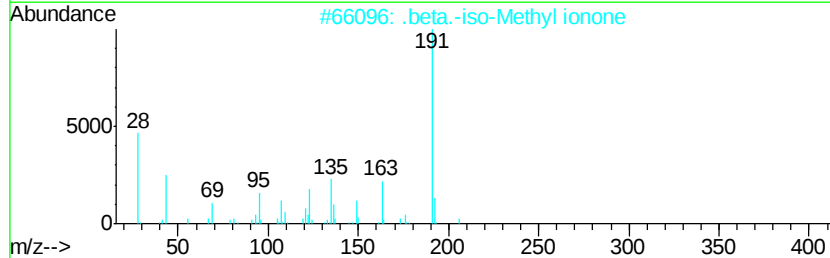
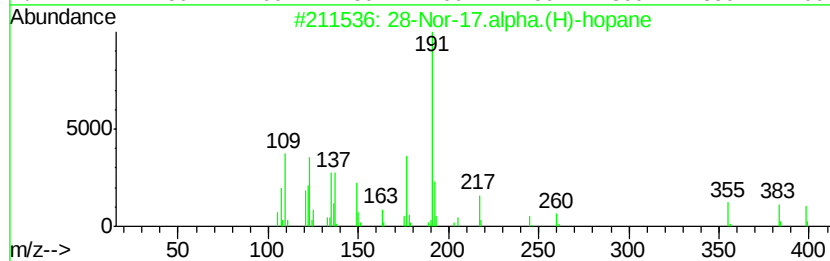
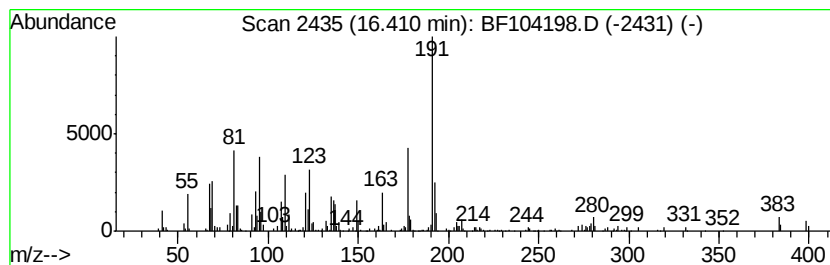
Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032818.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 28-Nor-17.alpha.(H)-hopane Concentration Rank 8

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.         |      |      |
|---------|-------------------------------------|--------------|------------------|--------------|------|------|
| 16.41   | 11.99 ng                            | 304078       | Perylene-d12     | 15.70        |      |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm      | CAS# | Qual |
| 1       | 28-Nor-17.alpha.(H)-hopane          | 398          | C29H50           | 053584-60-4  | 72   |      |
| 2       | .beta.-iso-Methyl ionone            | 206          | C14H22O          | 1000285-40-2 | 62   |      |
| 3       | 28-Nor-17.beta.(H)-hopane           | 398          | C29H50           | 036728-72-0  | 46   |      |
| 4       | 1-Cyclohexene, 1,3,3-trimethyl-2... | 206          | C14H22O          | 1000197-08-4 | 32   |      |
| 5       | Pentanoic acid, 4-methyl-2-(2,5-... | 278          | C13H18N4O3       | 1000294-63-5 | 30   |      |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

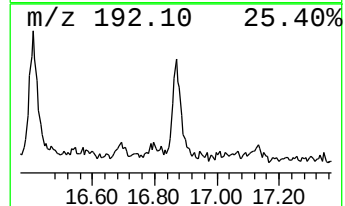
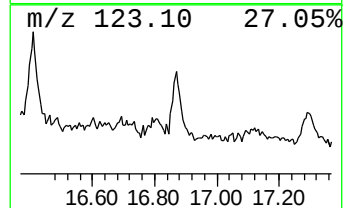
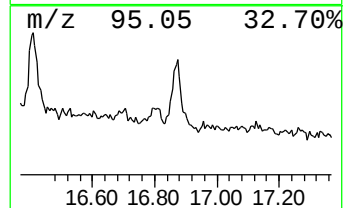
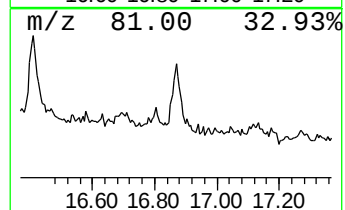
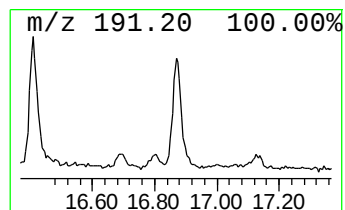
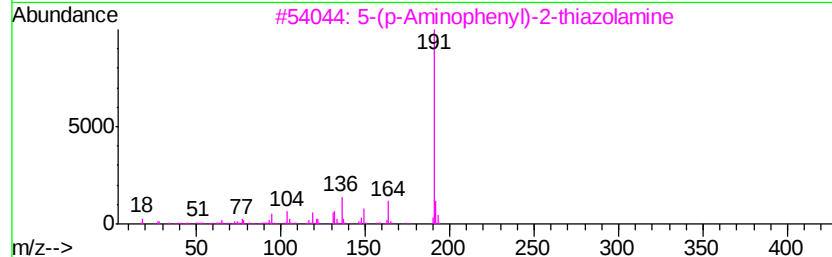
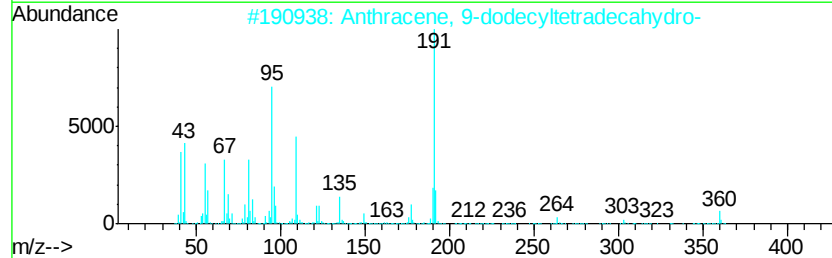
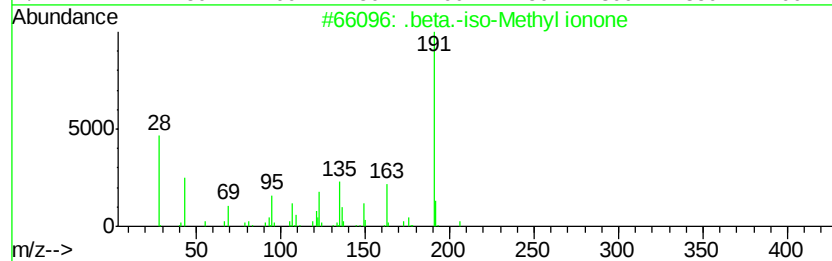
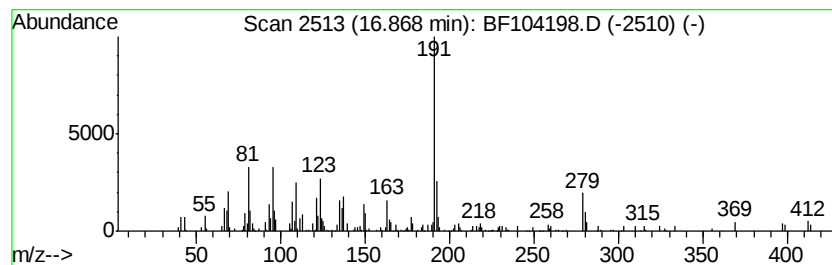
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown16.87 Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.87 | 7.47 ng | 189427 | Perylene-d12     | 15.70 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | .beta.-iso-Methyl ionone            | 206 | C14H22O   | 1000285-40-2 | 47   |
| 2    |    | Anthracene, 9-dodecyltetradecahy... | 360 | C26H48    | 055401-75-7  | 45   |
| 3    |    | 5-(p-Aminophenyl)-2-thiazolamine    | 191 | C9H9N3S   | 090349-87-4  | 38   |
| 4    |    | Isophthalic acid, 2-methyloct-5-... | 330 | C20H26O4  | 1000343-92-3 | 30   |
| 5    |    | 2-Fluoro-6-trifluoromethylbenzoi... | 320 | C14H6F6O2 | 1000357-69-2 | 27   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF040318\  
Data File : BF104198.D  
Acq On : 3 Apr 2018 23:42  
Operator : JU/SJ  
Sample : J2182-17  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

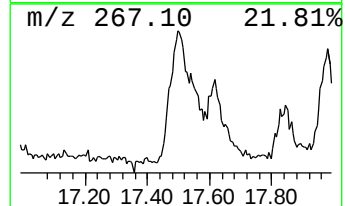
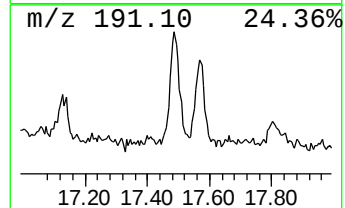
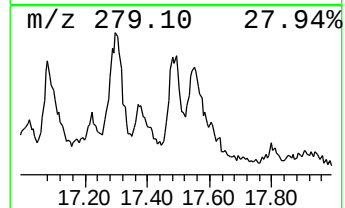
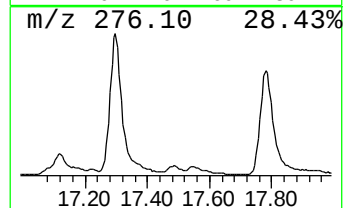
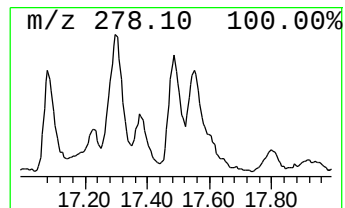
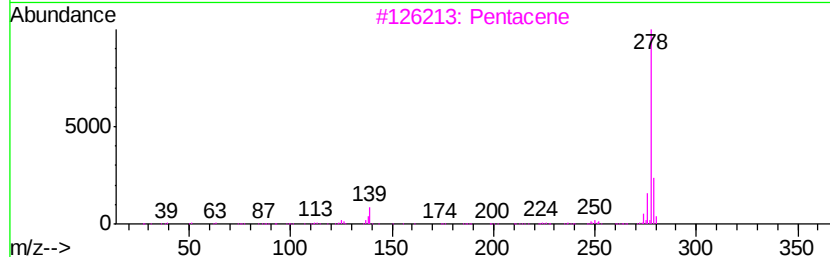
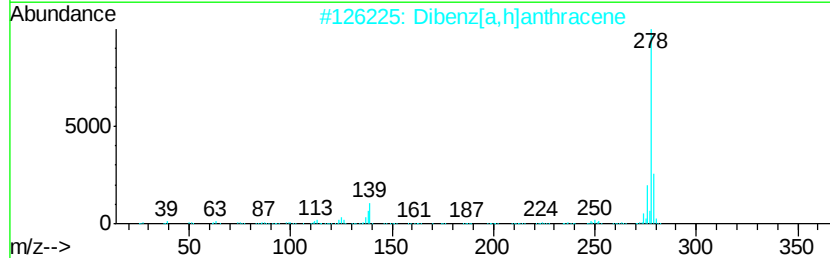
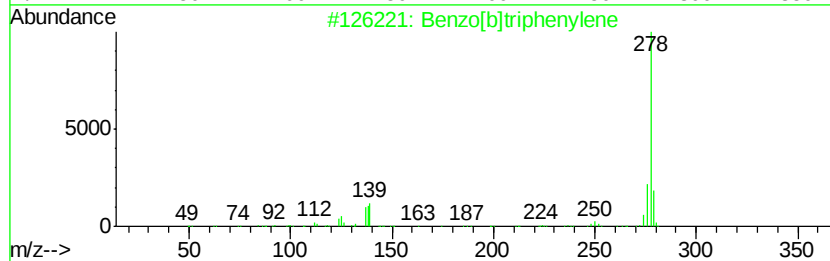
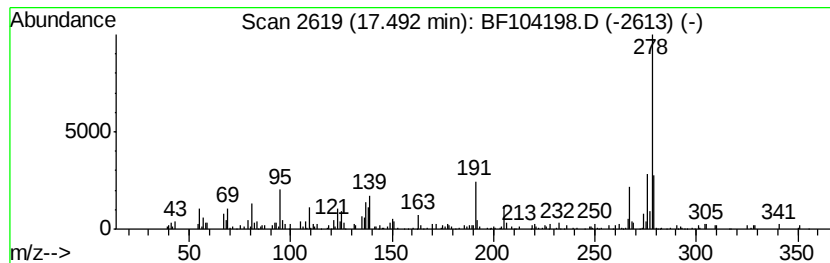
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Benzo[b]triphenylene Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.49 | 9.45 ng | 239507 | Perylene-d12     | 15.70 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[b]triphenylene  | 278 | C22H14  | 000215-58-7 | 91   |
| 2    |    | Dibenz[a,h]anthracene | 278 | C22H14  | 000053-70-3 | 90   |
| 3    |    | Pentacene             | 278 | C22H14  | 000135-48-8 | 78   |
| 4    |    | Benzo[b]chrysene      | 278 | C22H14  | 000214-17-5 | 78   |
| 5    |    | Picene                | 278 | C22H14  | 000213-46-7 | 78   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF040318\  
 Data File : BF104198.D  
 Acq On : 3 Apr 2018 23:42  
 Operator : JU/SJ  
 Sample : J2182-17  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 6-WM-DIP-7-WW(1-5)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF032818.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| unknown6.73          | 6.73  | 86.9    | ng    | 2381410  | 1                     | 6.96  | 547929  | 20.0 |
| Naphthalene, 1-me... | 9.05  | 8.1     | ng    | 290097   | 2                     | 8.24  | 714561  | 20.0 |
| unknown10.63         | 10.63 | 7.3     | ng    | 315401   | 3                     | 10.00 | 864158  | 20.0 |
| Dodecane             | 10.88 | 6.7     | ng    | 230049   | 4                     | 11.50 | 690861  | 20.0 |
| 9H-Fluorene, 2-me... | 11.10 | 6.5     | ng    | 226049   | 4                     | 11.50 | 690861  | 20.0 |
| Anthracene, 1,2,3... | 11.35 | 18.1    | ng    | 625745   | 4                     | 11.50 | 690861  | 20.0 |
| Dibenzothiophene     | 11.39 | 10.6    | ng    | 364806   | 4                     | 11.50 | 690861  | 20.0 |
| Phenanthrene, 2-m... | 12.00 | 16.6    | ng    | 575151   | 4                     | 11.50 | 690861  | 20.0 |
| 4H-Cyclopenta[def... | 12.12 | 46.0    | ng    | 1589060  | 4                     | 11.50 | 690861  | 20.0 |
| Naphthalene, 2-ph... | 12.29 | 11.5    | ng    | 397023   | 4                     | 11.50 | 690861  | 20.0 |
| Phenanthrene, 2,5... | 12.55 | 7.6     | ng    | 261476   | 4                     | 11.50 | 690861  | 20.0 |
| 11H-Benzo[b]fluorene | 13.28 | 8.8     | ng    | 1261450  | 5                     | 14.16 | 2868830 | 20.0 |
| Benzo[e]pyrene       | 15.36 | 8.2     | ng    | 208889   | 6                     | 15.70 | 507047  | 20.0 |
| 28-Nor-17.alpha.(... | 16.41 | 12.0    | ng    | 304078   | 6                     | 15.70 | 507047  | 20.0 |
| unknown16.87         | 16.87 | 7.5     | ng    | 189427   | 6                     | 15.70 | 507047  | 20.0 |
| Benzo[b]triphenylene | 17.49 | 9.4     | ng    | 239507   | 6                     | 15.70 | 507047  | 20.0 |