

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF022120\  
 Data File : BF119012.D  
 Acq On : 22 Feb 2020 00:37  
 Operator : CG/JU  
 Sample : L1613-01 5X  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-14-EP1-2

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF022020.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         | 6.392       | 729           | 732         | 743          | rVB      | 567545         | 596346        | 2.61%           | 0.180%        |
| 2         | 6.757       | 790           | 794         | 801          | rBV      | 895009         | 801620        | 3.50%           | 0.242%        |
| 3         | 7.016       | 834           | 838         | 846          | rBV      | 579936         | 634566        | 2.77%           | 0.192%        |
| 4         | 8.039       | 1007          | 1012        | 1013         | rVV      | 1062305        | 1033544       | 4.52%           | 0.312%        |
| 5         | 8.063       | 1013          | 1016        | 1019         | rVV      | 4029856        | 3392715       | 14.83%          | 1.025%        |
| 6         | 8.751       | 1129          | 1133        | 1136         | rBV      | 1989308        | 1561991       | 6.83%           | 0.472%        |
| 7         | 8.851       | 1145          | 1150        | 1153         | rVB      | 1180845        | 1098421       | 4.80%           | 0.332%        |
| 8         | 9.110       | 1191          | 1194        | 1197         | rVB      | 834031         | 666936        | 2.91%           | 0.201%        |
| 9         | 9.210       | 1205          | 1211        | 1215         | rBV2     | 620359         | 701567        | 3.07%           | 0.212%        |
| 10        | 9.374       | 1234          | 1239        | 1243         | rBV      | 606442         | 816490        | 3.57%           | 0.247%        |
| 11        | 9.451       | 1248          | 1252        | 1254         | rVV      | 1005367        | 968031        | 4.23%           | 0.292%        |
| 12        | 9.516       | 1259          | 1263        | 1268         | rVB      | 1059097        | 974883        | 4.26%           | 0.294%        |
| 13        | 9.663       | 1282          | 1288        | 1291         | rBV      | 6254970        | 7105655       | 31.06%          | 2.146%        |
| 14        | 9.786       | 1307          | 1309        | 1312         | rVV      | 1016950        | 902685        | 3.95%           | 0.273%        |
| 15        | 9.822       | 1312          | 1315        | 1318         | rVB2     | 743999         | 623607        | 2.73%           | 0.188%        |
| 16        | 9.992       | 1340          | 1344        | 1348         | rVV      | 2045386        | 1912865       | 8.36%           | 0.578%        |
| 17        | 10.104      | 1359          | 1363        | 1366         | rBV2     | 511464         | 673340        | 2.94%           | 0.203%        |
| 18        | 10.245      | 1384          | 1387        | 1389         | rVB      | 824205         | 725995        | 3.17%           | 0.219%        |
| 19        | 10.339      | 1400          | 1403        | 1409         | rVB      | 3357377        | 3427888       | 14.98%          | 1.035%        |
| 20        | 10.445      | 1418          | 1421        | 1426         | rVV      | 680273         | 685176        | 2.99%           | 0.207%        |
| 21        | 10.492      | 1426          | 1429        | 1432         | rVV      | 926996         | 919181        | 4.02%           | 0.278%        |
| 22        | 10.574      | 1437          | 1443        | 1448         | rBV2     | 1241662        | 1537076       | 6.72%           | 0.464%        |
| 23        | 10.704      | 1461          | 1465        | 1471         | rBV2     | 1191429        | 1298585       | 5.68%           | 0.392%        |
| 24        | 10.886      | 1491          | 1496        | 1499         | rBV2     | 800709         | 1101778       | 4.82%           | 0.333%        |
| 25        | 11.039      | 1519          | 1522        | 1527         | rVV      | 809624         | 1118104       | 4.89%           | 0.338%        |
| 26        | 11.086      | 1527          | 1530        | 1534         | rVV      | 1291661        | 1391955       | 6.08%           | 0.420%        |
| 27        | 11.121      | 1534          | 1536        | 1539         | rVV      | 522481         | 661242        | 2.89%           | 0.200%        |
| 28        | 11.168      | 1542          | 1544        | 1551         | rVB      | 1472284        | 1715388       | 7.50%           | 0.518%        |
| 29        | 11.321      | 1559          | 1570        | 1572         | rBV      | 8354884        | 13931264      | 60.89%          | 4.208%        |
| 30        | 11.363      | 1572          | 1577        | 1580         | rVV      | 5940456        | 6755767       | 29.53%          | 2.041%        |
| 31        | 11.386      | 1580          | 1581        | 1584         | rVB      | 905243         | 657920        | 2.88%           | 0.199%        |
| 32        | 11.504      | 1596          | 1601        | 1604         | rBV      | 1140390        | 1322596       | 5.78%           | 0.400%        |
| 33        | 11.568      | 1610          | 1612        | 1616         | rVB2     | 1096073        | 1014765       | 4.44%           | 0.307%        |
| 34        | 11.610      | 1616          | 1619        | 1623         | rVB3     | 991617         | 1008476       | 4.41%           | 0.305%        |

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Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-14-EP1-2

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

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF022020.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |          |        |        |
|----|--------|------|------|------|------|---------|----------|--------|--------|
| 35 | 11.692 | 1630 | 1633 | 1637 | rVB  | 890510  | 992467   | 4.34%  | 0.300% |
| 36 | 11.780 | 1644 | 1648 | 1651 | rVV  | 2574194 | 2765358  | 12.09% | 0.835% |
| 37 | 11.810 | 1651 | 1653 | 1656 | rVV  | 3445271 | 2922045  | 12.77% | 0.883% |
| 38 | 11.857 | 1658 | 1661 | 1664 | rVV  | 1557218 | 1615018  | 7.06%  | 0.488% |
| 39 | 11.898 | 1664 | 1668 | 1670 | rVV  | 4705339 | 5529774  | 24.17% | 1.670% |
| 40 | 11.916 | 1670 | 1671 | 1675 | rVB  | 2190239 | 1537239  | 6.72%  | 0.464% |
| 41 | 11.980 | 1680 | 1682 | 1685 | rVB2 | 739870  | 609415   | 2.66%  | 0.184% |
| 42 | 12.074 | 1694 | 1698 | 1700 | rVV  | 2245747 | 2175654  | 9.51%  | 0.657% |
| 43 | 12.110 | 1700 | 1704 | 1708 | rVB2 | 1201730 | 1502409  | 6.57%  | 0.454% |
| 44 | 12.257 | 1726 | 1729 | 1731 | rBV  | 691001  | 697113   | 3.05%  | 0.211% |
| 45 | 12.333 | 1738 | 1742 | 1745 | rBV  | 1986450 | 2521828  | 11.02% | 0.762% |
| 46 | 12.374 | 1745 | 1749 | 1755 | rVV3 | 1448760 | 2881164  | 12.59% | 0.870% |
| 47 | 12.439 | 1755 | 1760 | 1764 | rVB2 | 1740231 | 2995113  | 13.09% | 0.905% |
| 48 | 12.527 | 1764 | 1775 | 1777 | rBV  | 9594641 | 19433509 | 84.93% | 5.870% |
| 49 | 12.598 | 1784 | 1787 | 1792 | rVB  | 3953231 | 4415165  | 19.30% | 1.334% |
| 50 | 12.651 | 1792 | 1796 | 1798 | rBV  | 1677008 | 1906098  | 8.33%  | 0.576% |
| 51 | 12.757 | 1805 | 1814 | 1816 | rBV2 | 8928983 | 19924946 | 87.08% | 6.019% |
| 52 | 12.780 | 1816 | 1818 | 1822 | rVB2 | 1747698 | 1710989  | 7.48%  | 0.517% |
| 53 | 12.845 | 1826 | 1829 | 1834 | rBV2 | 1948614 | 2807818  | 12.27% | 0.848% |
| 54 | 12.886 | 1834 | 1836 | 1840 | rVB  | 1435211 | 1491661  | 6.52%  | 0.451% |
| 55 | 12.951 | 1841 | 1847 | 1853 | rBV4 | 2569169 | 5106075  | 22.32% | 1.542% |
| 56 | 13.062 | 1862 | 1866 | 1869 | rVB  | 4323406 | 6239573  | 27.27% | 1.885% |
| 57 | 13.098 | 1869 | 1872 | 1873 | rBV3 | 638060  | 619011   | 2.71%  | 0.187% |
| 58 | 13.127 | 1873 | 1877 | 1880 | rVV  | 3232536 | 3540240  | 15.47% | 1.069% |
| 59 | 13.168 | 1880 | 1884 | 1887 | rVV2 | 2951300 | 3772297  | 16.49% | 1.139% |
| 60 | 13.227 | 1890 | 1894 | 1895 | rBV2 | 1603015 | 1685421  | 7.37%  | 0.509% |
| 61 | 13.257 | 1897 | 1899 | 1902 | rVV  | 2130095 | 2303331  | 10.07% | 0.696% |
| 62 | 13.286 | 1902 | 1904 | 1907 | rVB  | 2212768 | 2066534  | 9.03%  | 0.624% |
| 63 | 13.457 | 1926 | 1933 | 1935 | rBV5 | 1313939 | 2371474  | 10.36% | 0.716% |
| 64 | 13.539 | 1944 | 1947 | 1948 | rBV  | 664051  | 697962   | 3.05%  | 0.211% |
| 65 | 13.568 | 1949 | 1952 | 1956 | rVB  | 1821727 | 2517505  | 11.00% | 0.760% |
| 66 | 13.680 | 1966 | 1971 | 1973 | rBV2 | 2329485 | 3537376  | 15.46% | 1.069% |
| 67 | 13.704 | 1973 | 1975 | 1978 | rVV  | 2726718 | 3014146  | 13.17% | 0.910% |
| 68 | 13.733 | 1978 | 1980 | 1986 | rVV3 | 2962154 | 4567874  | 19.96% | 1.380% |
| 69 | 13.792 | 1986 | 1990 | 1994 | rVB2 | 1657776 | 2708521  | 11.84% | 0.818% |
| 70 | 13.945 | 2007 | 2016 | 2018 | rBV2 | 7841034 | 19110171 | 83.52% | 5.773% |
| 71 | 13.986 | 2021 | 2023 | 2027 | rVV2 | 8065104 | 7271539  | 31.78% | 2.197% |

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 TP-14-EP1-2

Integration Parameters: rteint.p  
 Integrator: RTE  
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 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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF022020.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |          |         |        |
|-----|--------|------|------|------|------|---------|----------|---------|--------|
| 72  | 14.045 | 2031 | 2033 | 2037 | rVV3 | 1061291 | 1149310  | 5.02%   | 0.347% |
| 73  | 14.086 | 2037 | 2040 | 2045 | rVV3 | 2219957 | 3341596  | 14.60%  | 1.009% |
| 74  | 14.127 | 2045 | 2047 | 2049 | rVB  | 1104129 | 926142   | 4.05%   | 0.280% |
| 75  | 14.157 | 2049 | 2052 | 2056 | rBV2 | 1849636 | 2571671  | 11.24%  | 0.777% |
| 76  | 14.280 | 2071 | 2073 | 2075 | rBV2 | 896454  | 1020254  | 4.46%   | 0.308% |
| 77  | 14.321 | 2077 | 2080 | 2083 | rVB  | 2687066 | 2935007  | 12.83%  | 0.887% |
| 78  | 14.357 | 2083 | 2086 | 2088 | rVB2 | 1746465 | 1617627  | 7.07%   | 0.489% |
| 79  | 14.398 | 2089 | 2093 | 2095 | rBV3 | 883648  | 1329003  | 5.81%   | 0.401% |
| 80  | 14.421 | 2095 | 2097 | 2099 | rVB2 | 927185  | 746289   | 3.26%   | 0.225% |
| 81  | 14.451 | 2099 | 2102 | 2104 | rBV  | 2033119 | 2380915  | 10.41%  | 0.719% |
| 82  | 14.509 | 2109 | 2112 | 2115 | rVV2 | 1418317 | 1422568  | 6.22%   | 0.430% |
| 83  | 14.557 | 2118 | 2120 | 2127 | rVB2 | 549418  | 690201   | 3.02%   | 0.208% |
| 84  | 14.645 | 2130 | 2135 | 2141 | rVB4 | 795910  | 2006423  | 8.77%   | 0.606% |
| 85  | 14.809 | 2159 | 2163 | 2168 | rBV2 | 1108812 | 1847023  | 8.07%   | 0.558% |
| 86  | 14.998 | 2184 | 2195 | 2201 | rBV  | 6849014 | 22880774 | 100.00% | 6.912% |
| 87  | 15.092 | 2205 | 2211 | 2219 | rVB  | 3485905 | 5901521  | 25.79%  | 1.783% |
| 88  | 15.162 | 2219 | 2223 | 2227 | rBV2 | 762752  | 1540025  | 6.73%   | 0.465% |
| 89  | 15.280 | 2235 | 2243 | 2247 | rBV  | 3978862 | 9194870  | 40.19%  | 2.777% |
| 90  | 15.362 | 2247 | 2257 | 2260 | rVV2 | 6272599 | 14263127 | 62.34%  | 4.308% |
| 91  | 15.433 | 2264 | 2269 | 2273 | rVB2 | 3449681 | 5023675  | 21.96%  | 1.517% |
| 92  | 15.556 | 2285 | 2290 | 2293 | rBV  | 931604  | 1750594  | 7.65%   | 0.529% |
| 93  | 15.733 | 2316 | 2320 | 2325 | rVB2 | 755347  | 1150631  | 5.03%   | 0.348% |
| 94  | 15.927 | 2349 | 2353 | 2362 | rVB2 | 1123063 | 1977607  | 8.64%   | 0.597% |
| 95  | 16.845 | 2495 | 2509 | 2513 | rBV2 | 3233978 | 10998746 | 48.07%  | 3.322% |
| 96  | 16.880 | 2513 | 2515 | 2523 | rVB  | 801506  | 970240   | 4.24%   | 0.293% |
| 97  | 16.974 | 2525 | 2531 | 2536 | rBV  | 1128894 | 2534141  | 11.08%  | 0.765% |
| 98  | 17.033 | 2537 | 2541 | 2546 | rBV2 | 606184  | 1113758  | 4.87%   | 0.336% |
| 99  | 17.292 | 2571 | 2585 | 2592 | rVB  | 2701821 | 8817432  | 38.54%  | 2.663% |
| 100 | 17.497 | 2612 | 2620 | 2630 | rVB2 | 1296910 | 3646696  | 15.94%  | 1.102% |

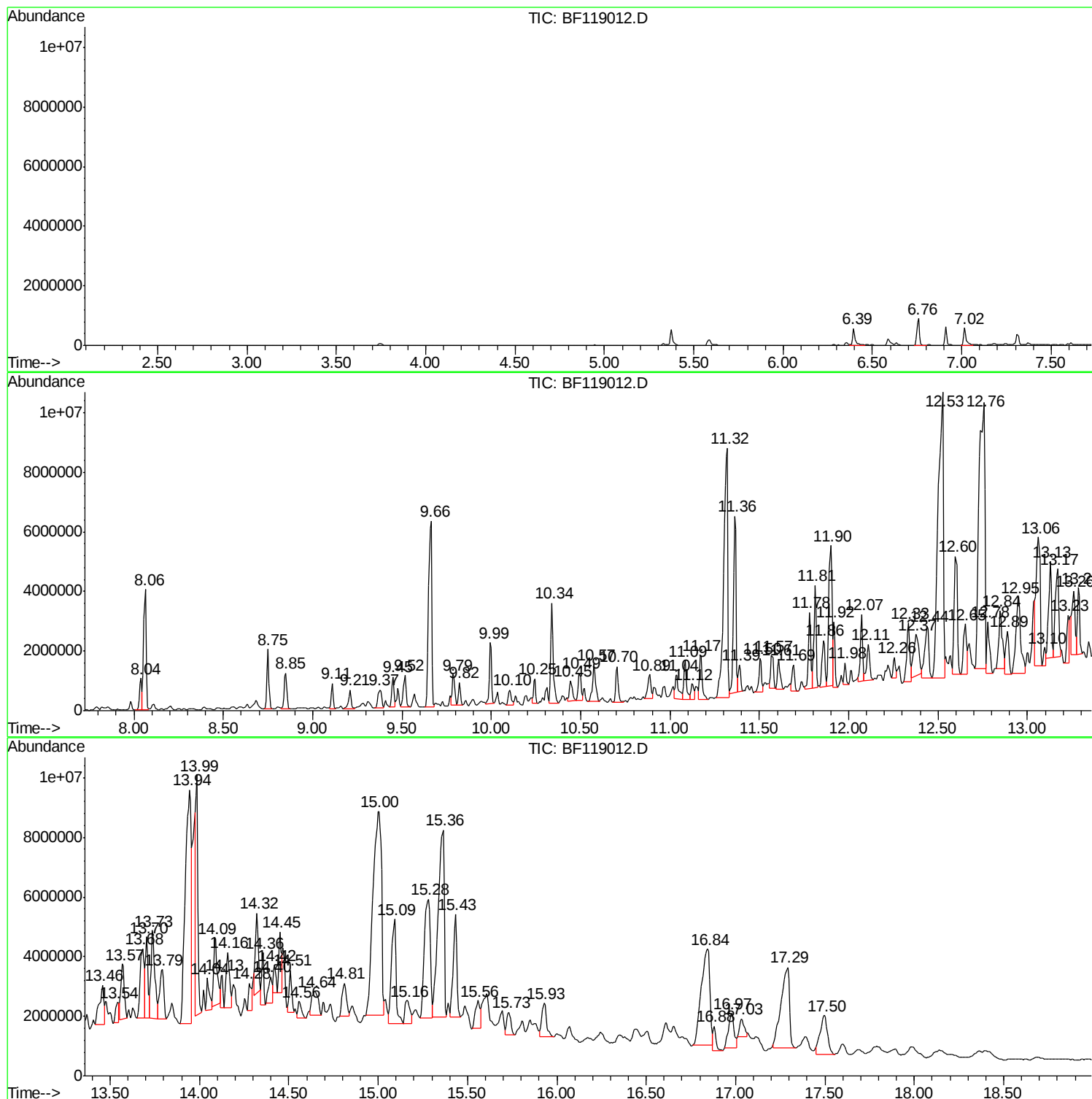
Sum of corrected areas: 331050117

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Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

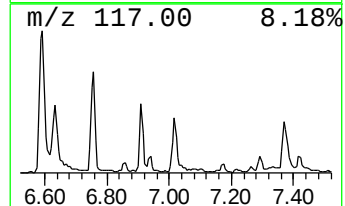
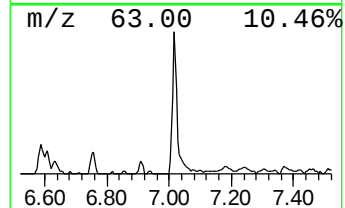
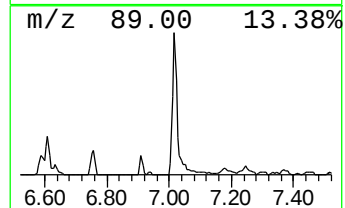
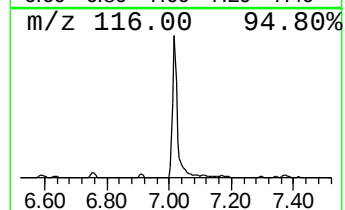
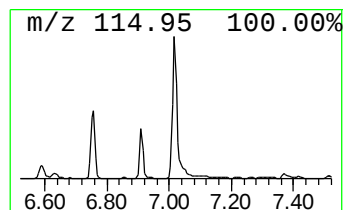
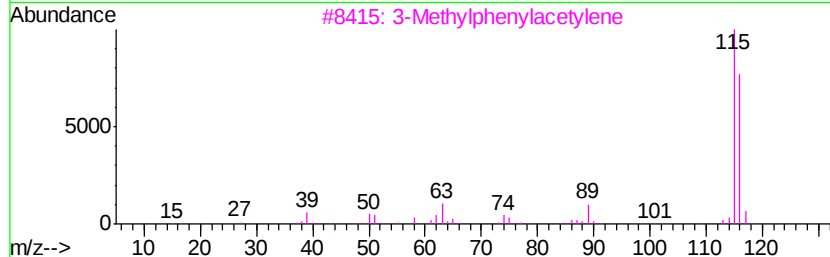
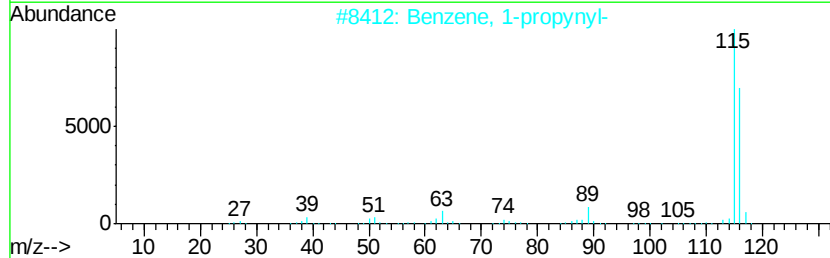
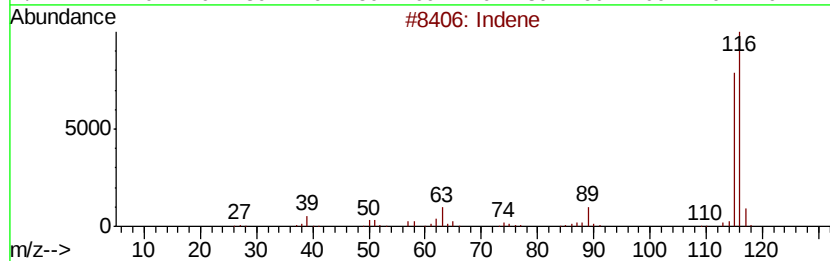
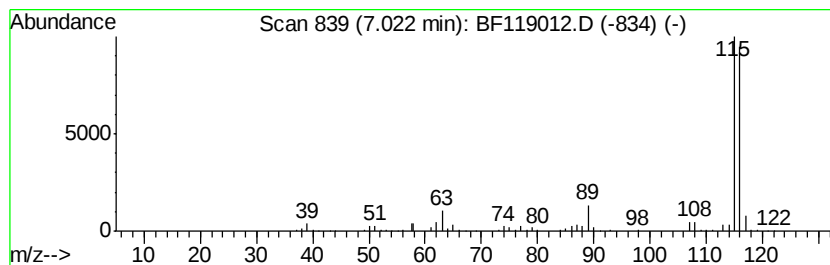
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 Indene Concentration Rank 6

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.02 | 15.83 ng | 634566 | 1,4-Dichlorobenzene-d4 | 6.76 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Indene                       | 116 | C9H8    | 000095-13-6 | 97   |
| 2    |    | Benzene, 1-propynyl-         | 116 | C9H8    | 000673-32-5 | 95   |
| 3    |    | 3-Methylphenylacetylene      | 116 | C9H8    | 000766-82-5 | 94   |
| 4    |    | Benzene, 1-ethynyl-4-methyl- | 116 | C9H8    | 000766-97-2 | 91   |
| 5    |    | Benzene, 1,2-propadienyl-    | 116 | C9H8    | 002327-99-3 | 91   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

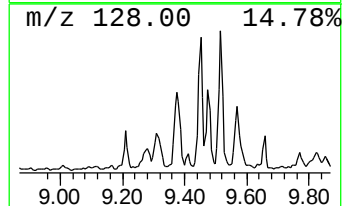
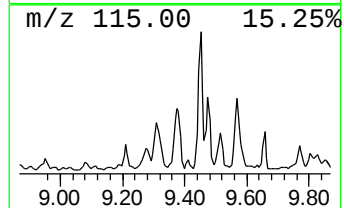
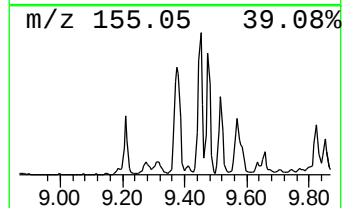
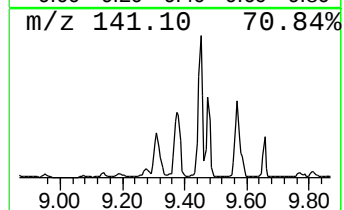
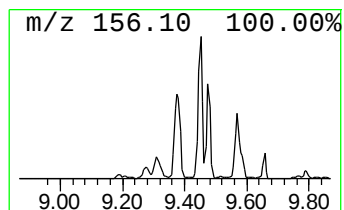
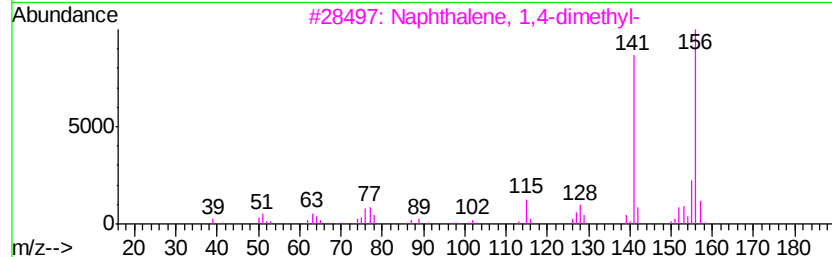
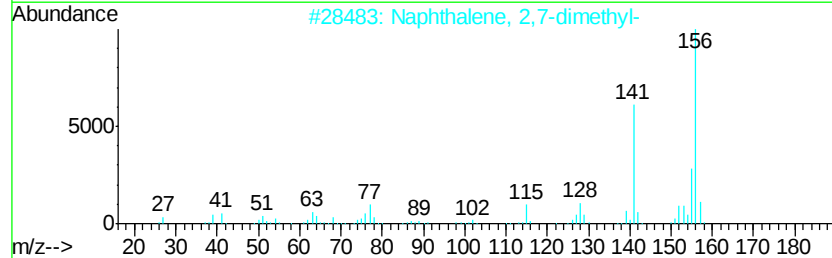
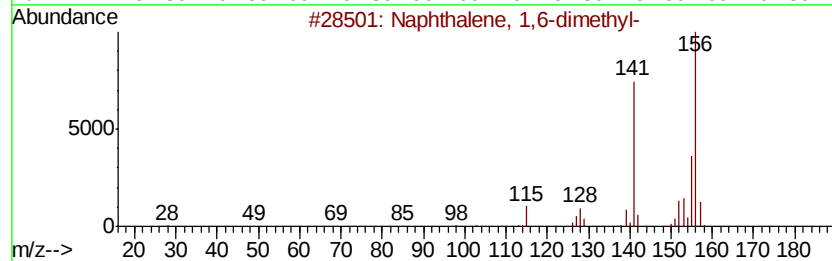
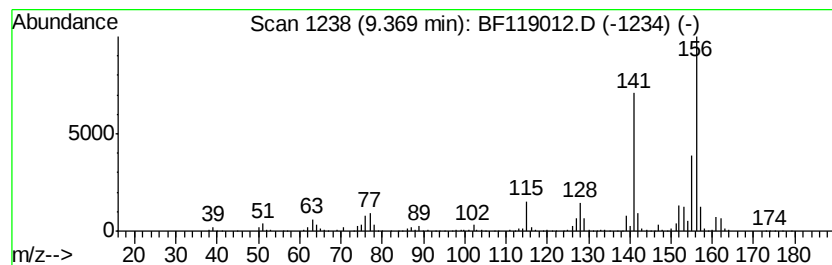
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Naphthalene, 1,6-dimethyl- Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.37 | 18.09 ng | 816490 | Acenaphthene-d10 | 9.79 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 2    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |
| 3    |    | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 96   |
| 4    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 95   |
| 5    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

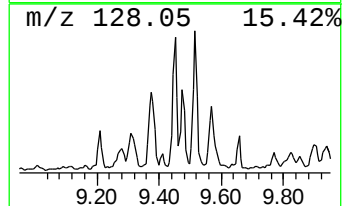
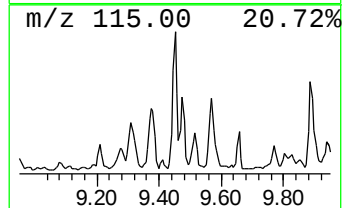
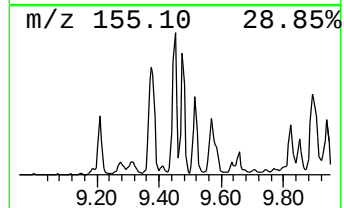
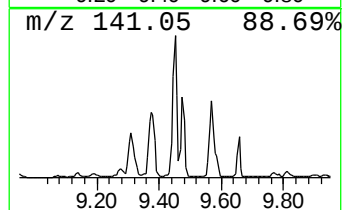
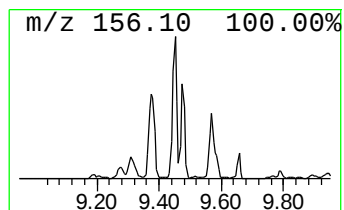
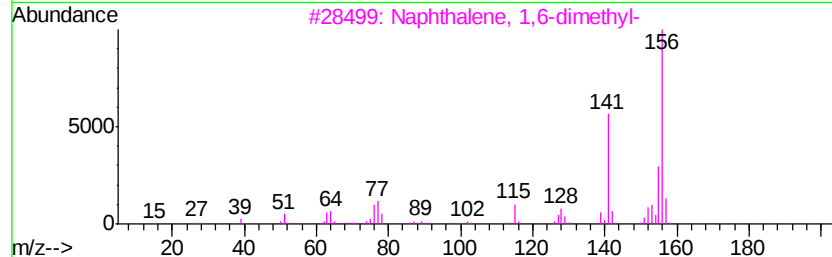
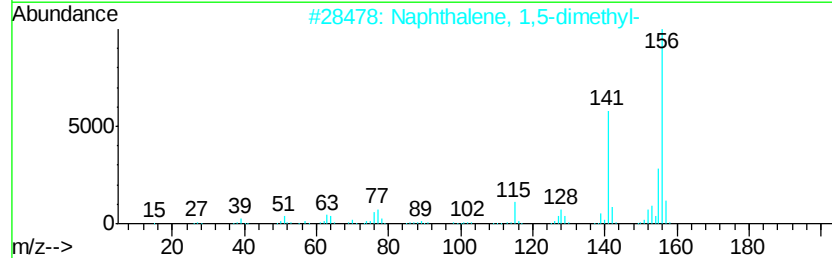
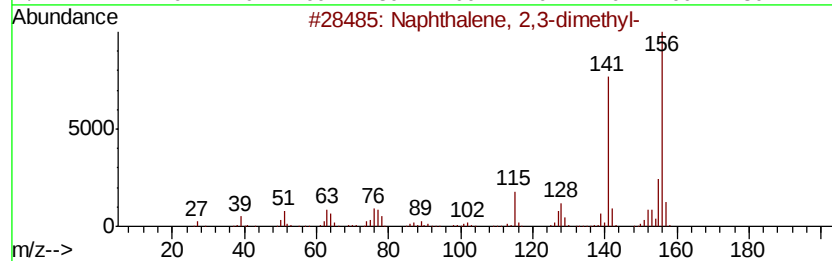
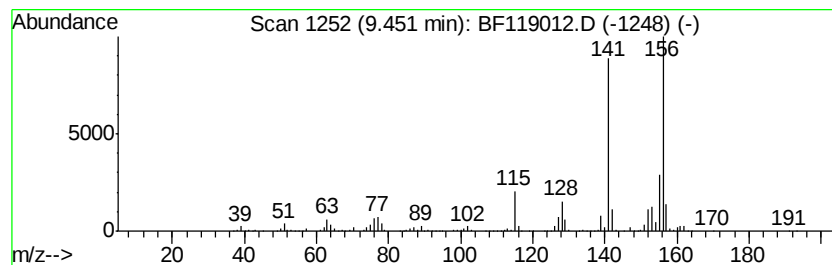
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Naphthalene, 2,3-dimethyl- Concentration Rank 2

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.45 | 21.45 ng | 968031 | Acenaphthene-d10 | 9.79 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 2    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |
| 3    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |
| 4    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 5    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

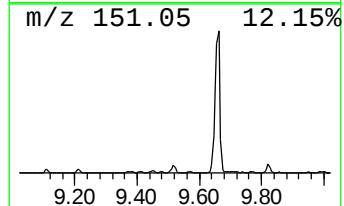
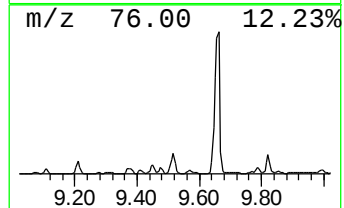
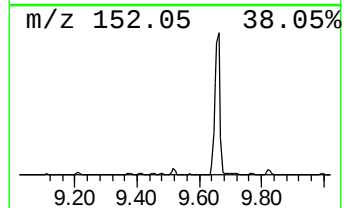
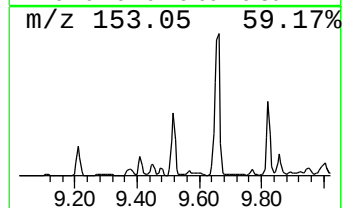
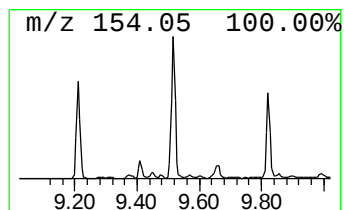
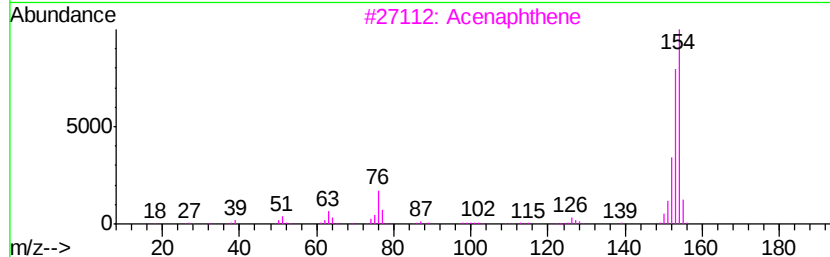
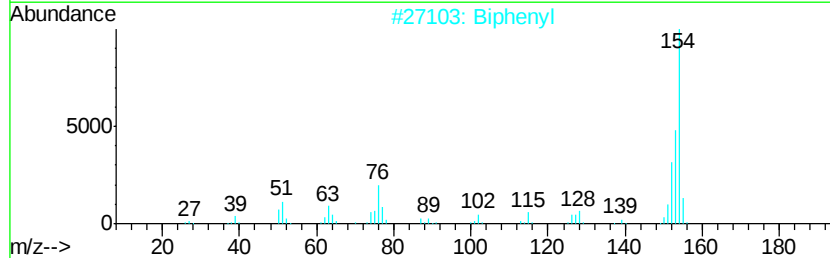
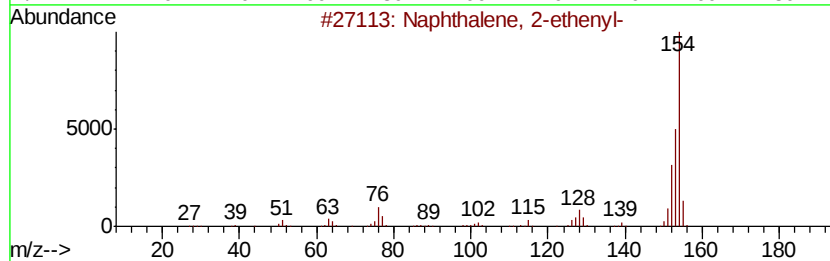
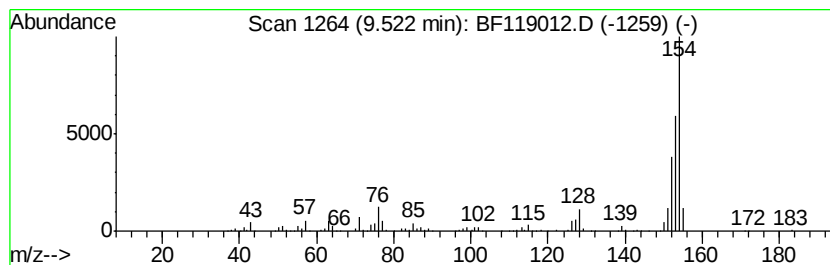
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Naphthalene, 2-ethenyl- Concentration Rank 1

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.52 | 21.60 ng | 974883 | Acenaphthene-d10 | 9.79 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Naphthalene, 2-ethenyl-             | 154 | C12H10  | 000827-54-3  | 96   |
| 2    |    | Biphenyl                            | 154 | C12H10  | 000092-52-4  | 90   |
| 3    |    | Acenaphthene                        | 154 | C12H10  | 000083-32-9  | 81   |
| 4    |    | 1,4-Ethenonaphthalene, 1,4-dihydro- | 154 | C12H10  | 007322-47-6  | 62   |
| 5    |    | 1-(2,2-Difluoroethenyl)-4-methyl... | 154 | C9H8F2  | 1000305-64-2 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

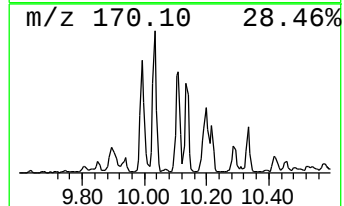
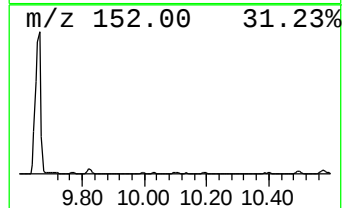
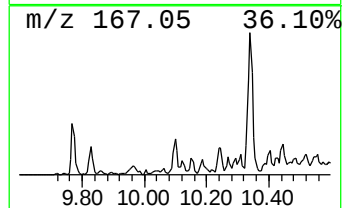
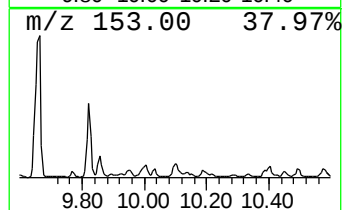
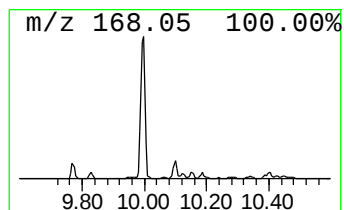
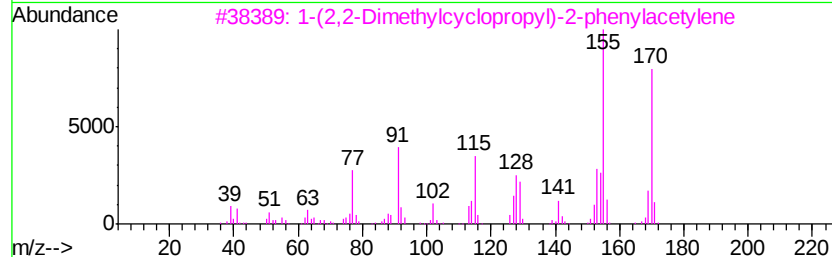
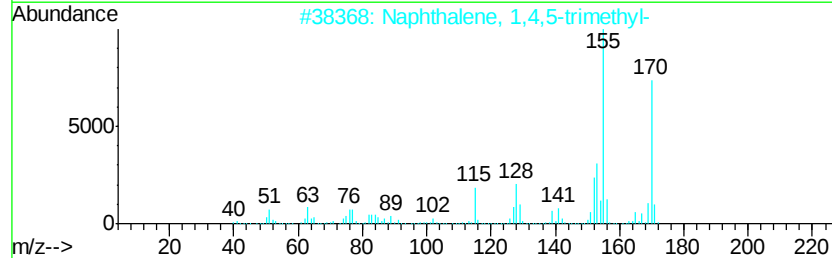
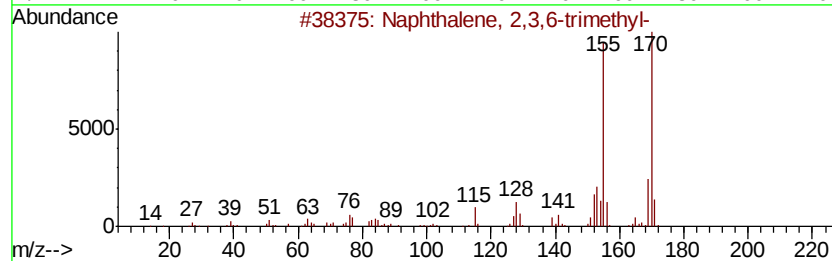
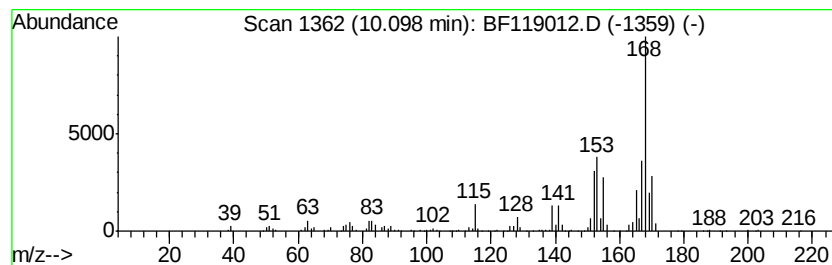
Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Naphthalene, 2,3,6-trimethyl- Concentration Rank 8

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.10     | 14.92 ng                            | 673340 | Acenaphthene-d10 | 9.79         |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Naphthalene, 2,3,6-trimethyl-       | 170    | C13H14           | 000829-26-5  | 91   |
| 2         | Naphthalene, 1,4,5-trimethyl-       | 170    | C13H14           | 002131-41-1  | 87   |
| 3         | 1-(2,2-Dimethylcyclopropyl)-2-ph... | 170    | C13H14           | 1000294-91-1 | 60   |
| 4         | Naphthalene, 1,6,7-trimethyl-       | 170    | C13H14           | 002245-38-7  | 55   |
| 5         | 3-(2-Methyl-propenyl)-1H-indene     | 170    | C13H14           | 1000187-78-5 | 55   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

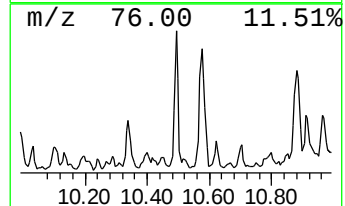
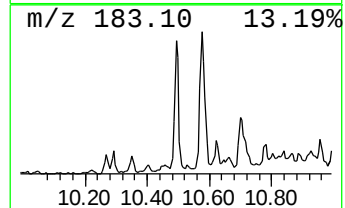
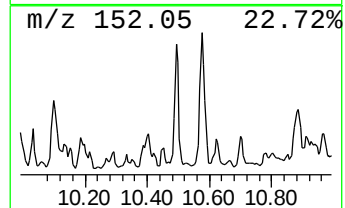
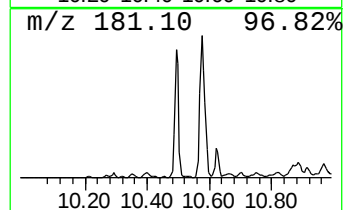
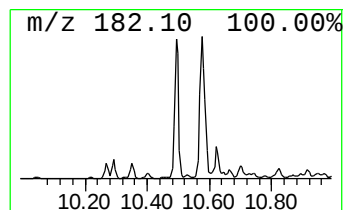
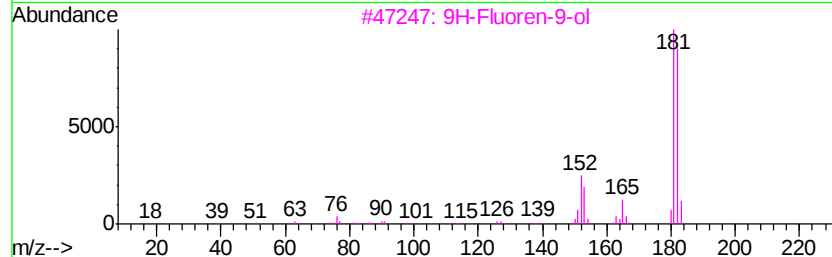
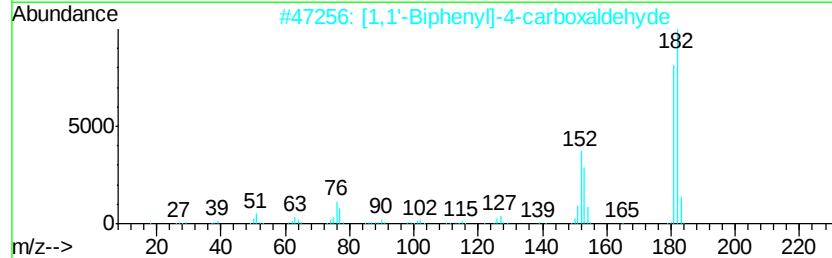
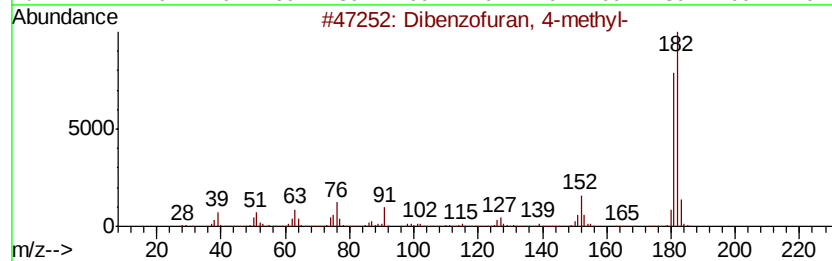
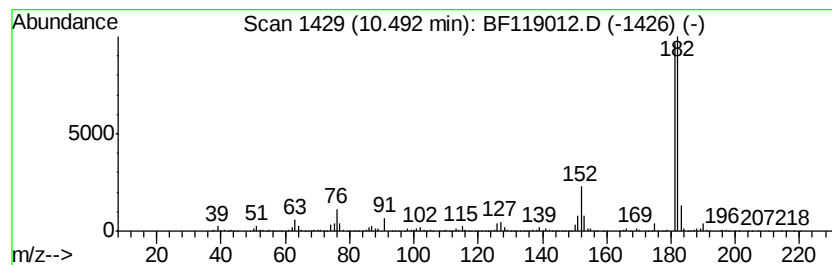
Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 8 Dibenzofuran, 4-methyl- Concentration Rank 3

| R.T.    | EstConc  | Area                             | Relative to ISTD | R.T.            |
|---------|----------|----------------------------------|------------------|-----------------|
| 10.49   | 20.37 ng | 919181                           | Acenaphthene-d10 | 9.79            |
| Hit# of | 5        | Tentative ID                     | MW MolForm       | CAS# Qual       |
| 1       |          | Dibenzofuran, 4-methyl-          | 182 C13H10O      | 007320-53-8 93  |
| 2       |          | [1,1'-Biphenyl]-4-carboxaldehyde | 182 C13H10O      | 003218-36-8 90  |
| 3       |          | 9H-Fluoren-9-ol                  | 182 C13H10O      | 001689-64-1 80  |
| 4       |          | 3-(2-Naphthyl)acrylaldehyde      | 182 C13H10O      | 020884-05-3 72  |
| 5       |          | Norharmene, N-methyl-            | 182 C12H10N2     | 1000374-78-6 64 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

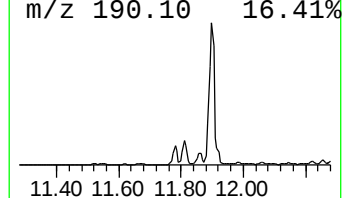
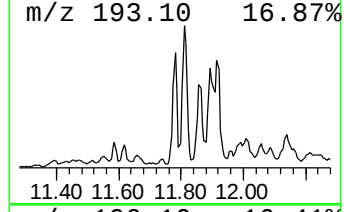
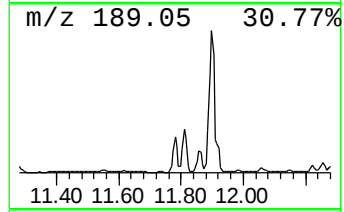
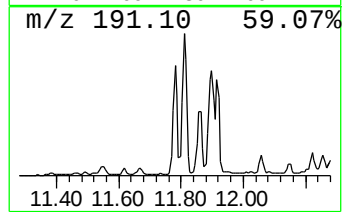
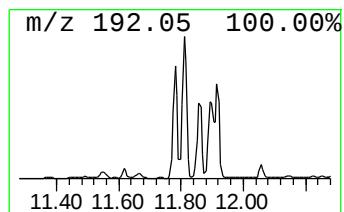
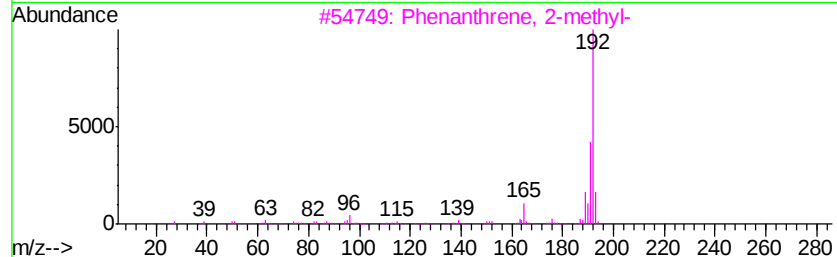
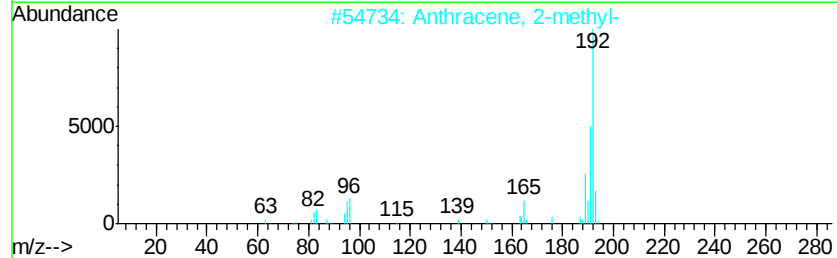
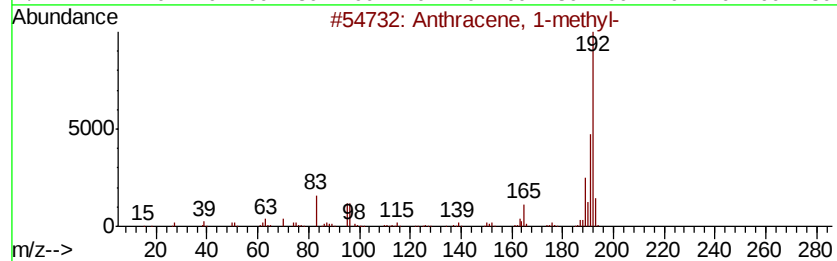
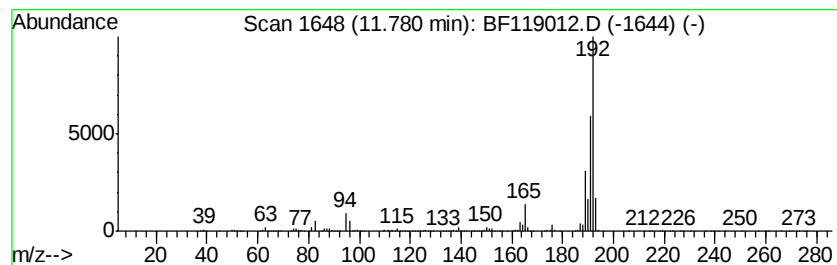
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Anthracene, 1-methyl- Concentration Rank 20

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 11.78 | 3.97 ng | 2765360 | Phenanthrene-d10 | 11.28 |

| Hit# | of | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 1-methyl-                 | 192 | C15H12  | 000610-48-0 | 94   |
| 2    |    | Anthracene, 2-methyl-                 | 192 | C15H12  | 000613-12-7 | 93   |
| 3    |    | Phenanthrene, 2-methyl-               | 192 | C15H12  | 002531-84-2 | 92   |
| 4    |    | Phenanthrene, 1-methyl-               | 192 | C15H12  | 000832-69-9 | 91   |
| 5    |    | 1H-Cyclopropa[1]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

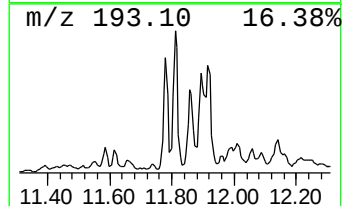
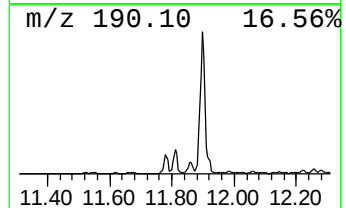
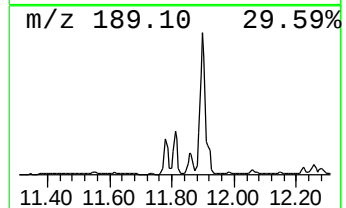
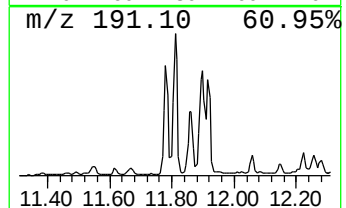
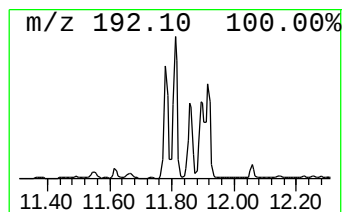
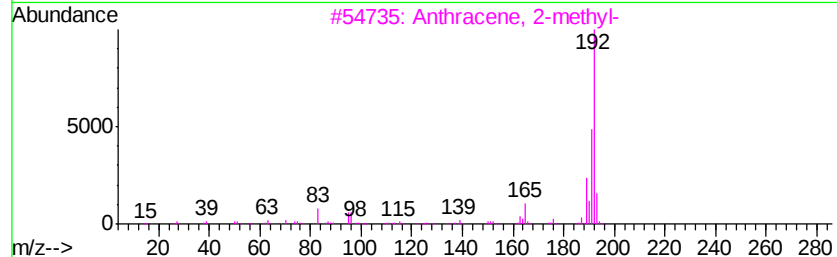
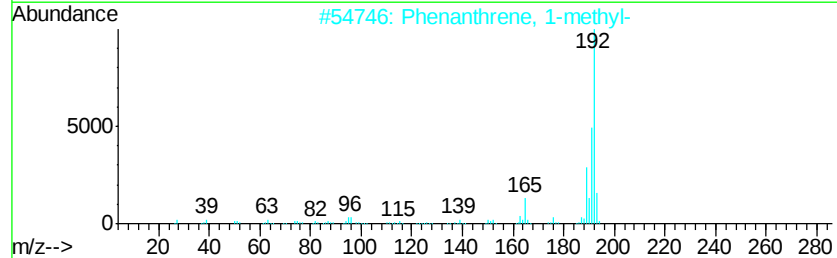
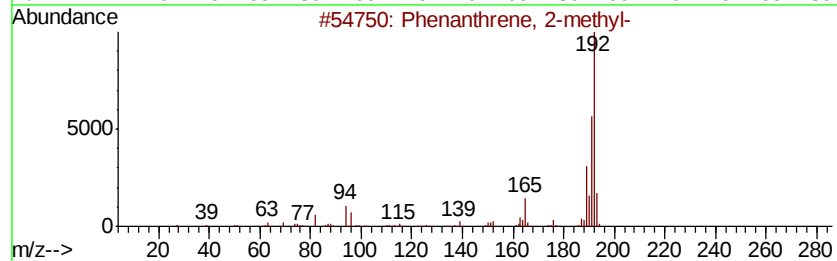
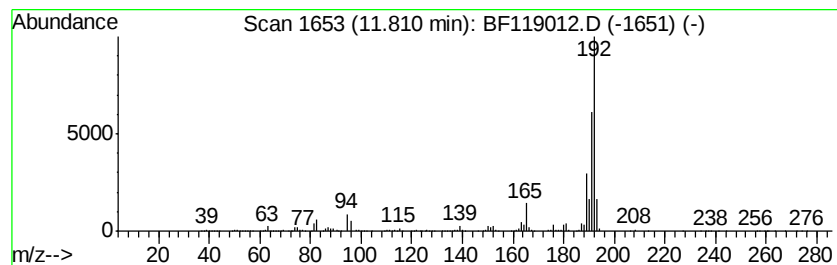
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Phenanthrene, 2-methyl- Concentration Rank 18

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 11.81 | 4.19 ng | 2922050 | Phenanthrene-d10 | 11.28 |

| Hit# | of | 5 | Tentative ID                           | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2-methyl-                | 192 | C15H12  | 002531-84-2 | 98   |
| 2    |    |   | Phenanthrene, 1-methyl-                | 192 | C15H12  | 000832-69-9 | 95   |
| 3    |    |   | Anthracene, 2-methyl-                  | 192 | C15H12  | 000613-12-7 | 94   |
| 4    |    |   | 1H-Cyclopropa[1,1]phenanthrene, 1a,... | 192 | C15H12  | 000949-41-7 | 93   |
| 5    |    |   | Anthracene, 1-methyl-                  | 192 | C15H12  | 000610-48-0 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

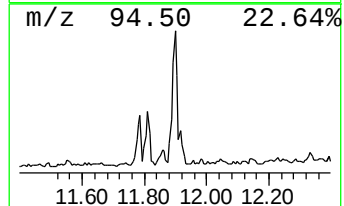
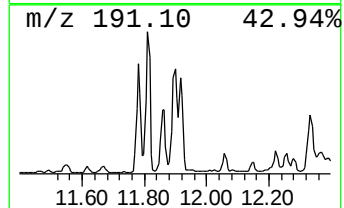
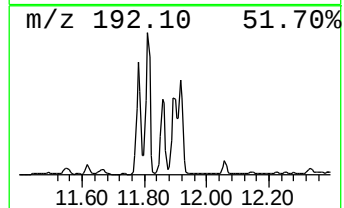
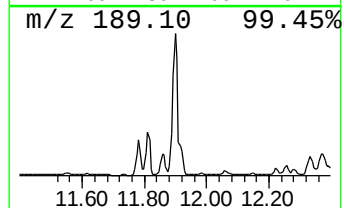
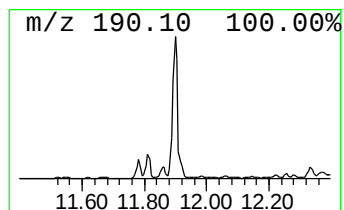
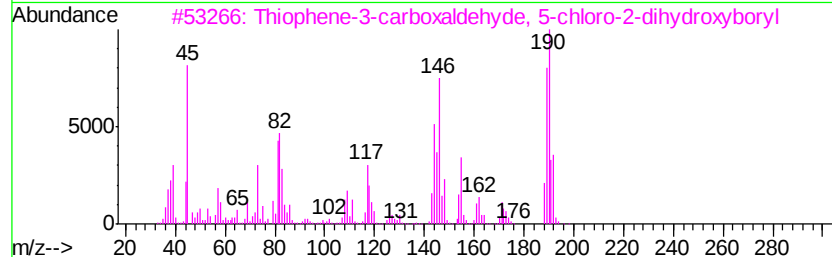
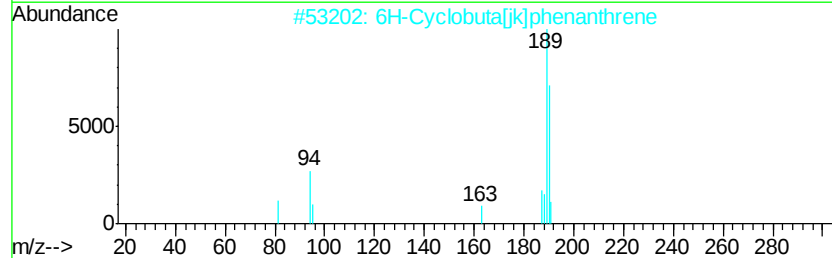
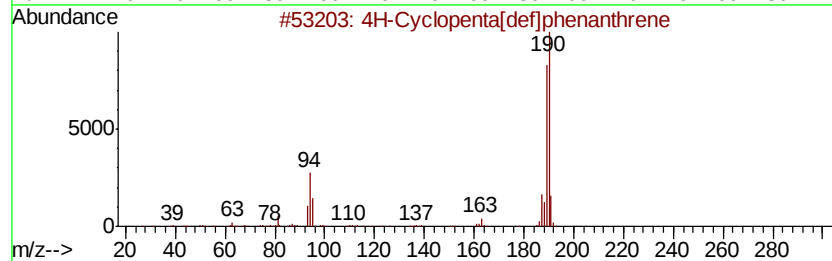
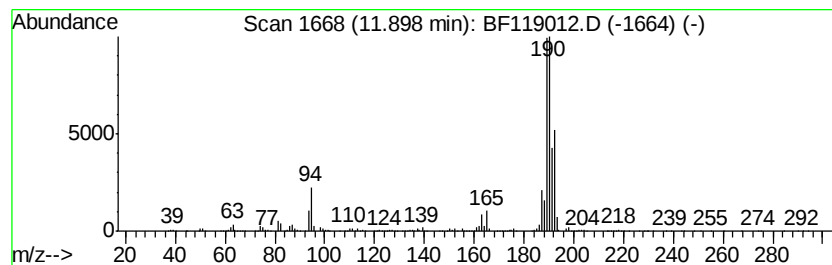
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ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 11

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 11.90     | 7.94 ng                             | 5529770 | Phenanthrene-d10 | 11.28       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | 4H-Cyclopenta[def]phenanthrene      | 190     | C15H10           | 000203-64-5 | 92   |
| 2         | 6H-Cyclobuta[ik]phenanthrene        | 190     | C15H10           | 083469-43-6 | 58   |
| 3         | Thiophene-3-carboxaldehyde, 5-ch... | 190     | C5H4BClO3S       | 036155-87-0 | 50   |
| 4         | 2,3,5,6-Tetramethylterephthalald... | 190     | C12H14O2         | 007072-01-7 | 43   |
| 5         | Methyl diselenide                   | 190     | C2H6Se2          | 007101-31-7 | 41   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

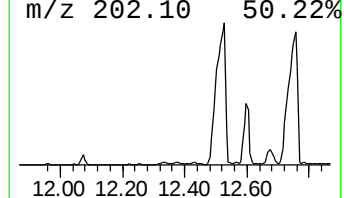
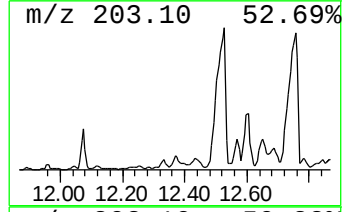
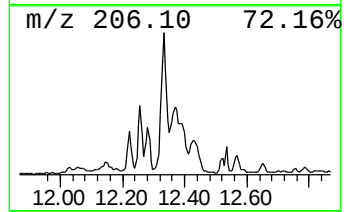
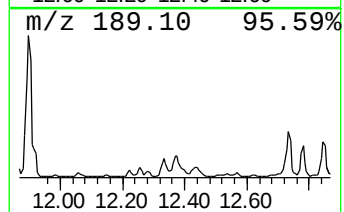
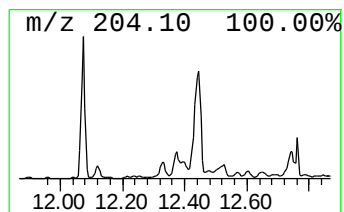
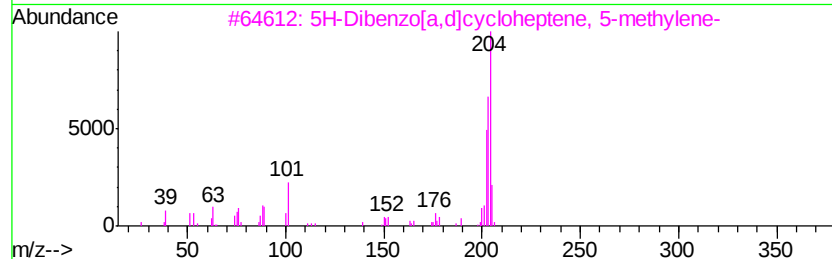
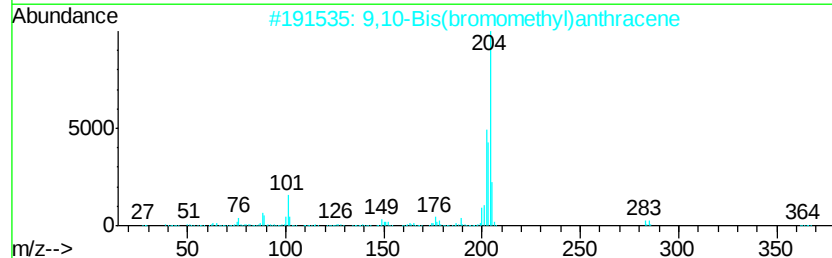
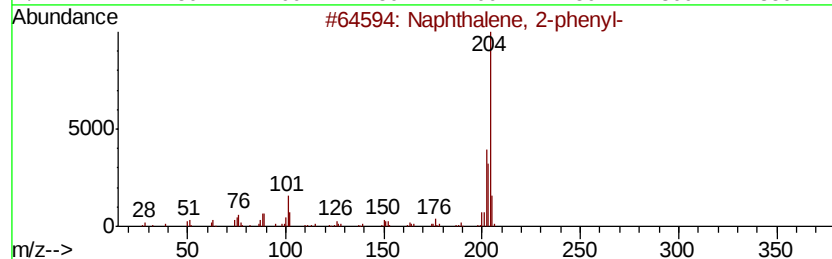
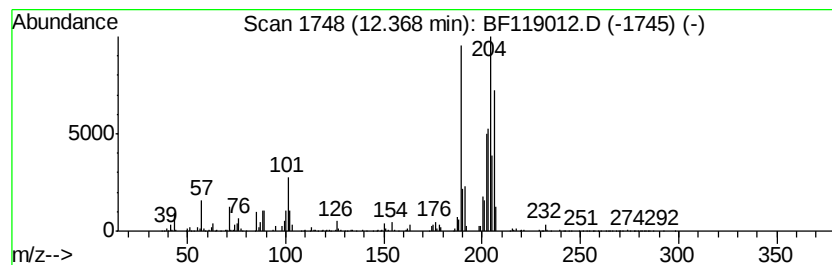
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 unknown12.37 Concentration Rank 19

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 12.37 | 4.14 ng | 2881160 | Phenanthrene-d10 | 11.28 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Naphthalene, 2-phenyl-              | 204 | C16H12    | 000612-94-2  | 42   |
| 2    |    | 9,10-Bis(bromomethyl)anthracene     | 362 | C16H12Br2 | 034373-96-1  | 38   |
| 3    |    | 5H-Dibenzo[a,d]cycloheptene, 5-m... | 204 | C16H12    | 002975-79-3  | 30   |
| 4    |    | Quinoline, 8-methoxy-5-nitro-       | 204 | C10H8N2O3 | 1000298-88-7 | 30   |
| 5    |    | Acephenanthrylene, 4,5-dihydro-     | 204 | C16H12    | 006232-48-0  | 25   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

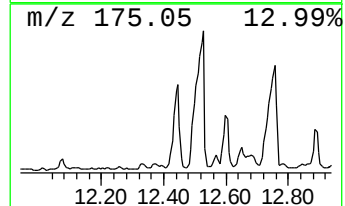
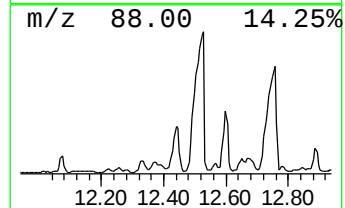
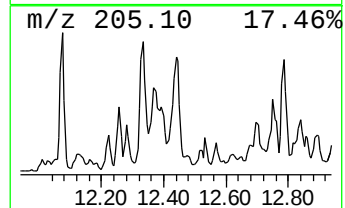
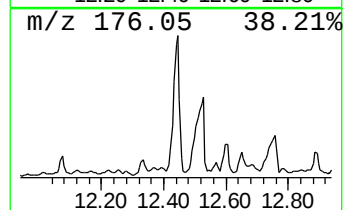
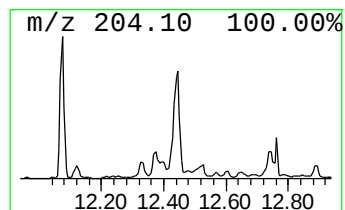
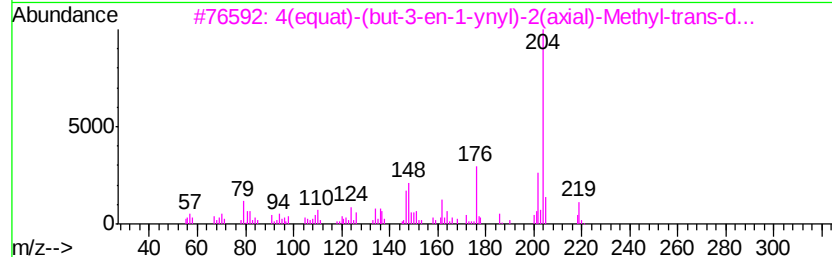
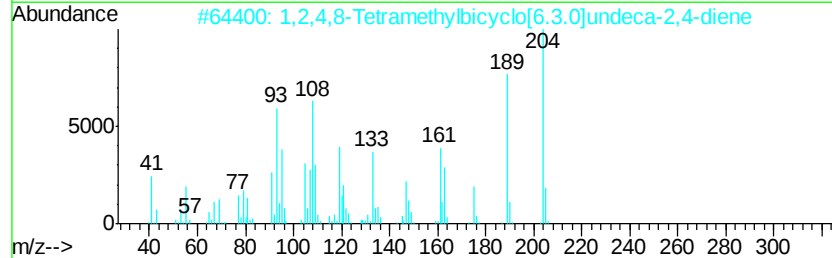
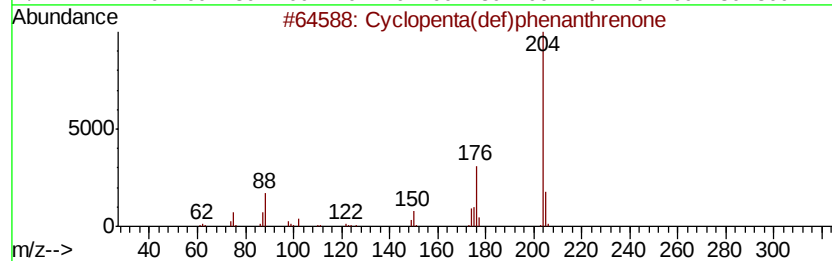
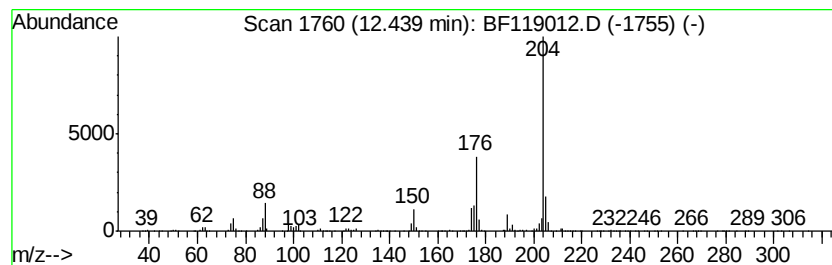
Instrument :  
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ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 Cyclopenta(def)phenanthrenone Concentration Rank 17

| R.T.    | EstConc | Area                                                                                  | Relative to ISTD | R.T.      |              |      |
|---------|---------|---------------------------------------------------------------------------------------|------------------|-----------|--------------|------|
| 12.44   | 4.30 ng | 2995110                                                                               | Phenanthrene-d10 | 11.28     |              |      |
| Hit# of | 5       | Tentative ID                                                                          | MW               | MolForm   | CAS#         | Qual |
| 1       |         | Cvclopenta(def)phenanthrenone                                                         | 204              | C15H8O    | 005737-13-3  | 97   |
| 2       |         | 1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene                                     | 204              | C15H24    | 137235-51-9  | 55   |
| 3       |         | 4(equat)-(but-3-en-1-ynyl)-2(axial)-Methyl-trans-dibenzylidene-1,2,3,4,5,6-hexatriene | 219              | C14H21NO  | 038423-22-2  | 50   |
| 4       |         | [1,1'-Biphenyl]-4,4'-dicarbonitrile                                                   | 204              | C14H8N2   | 001591-30-6  | 47   |
| 5       |         | (E)-Ethyl 3-oxo-2-(pyrid-2-yl)methylpentanoate                                        | 219              | C12H13NO3 | 1000147-59-7 | 45   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

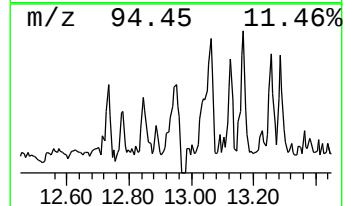
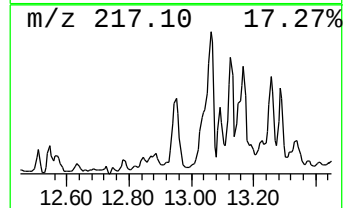
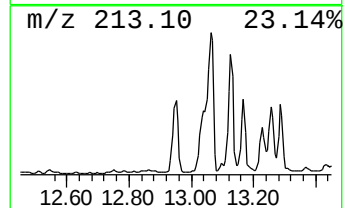
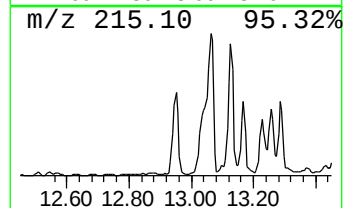
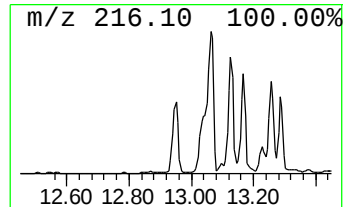
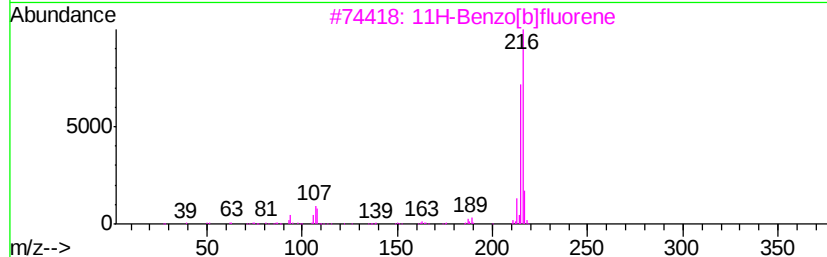
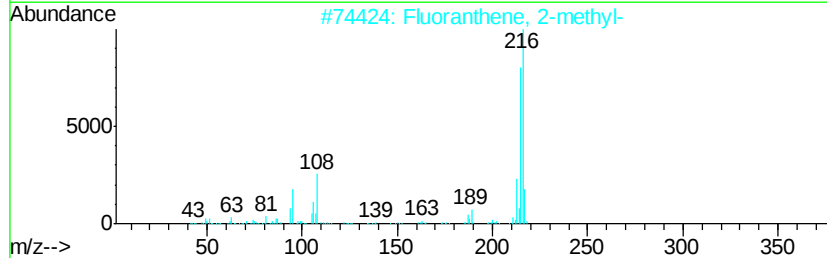
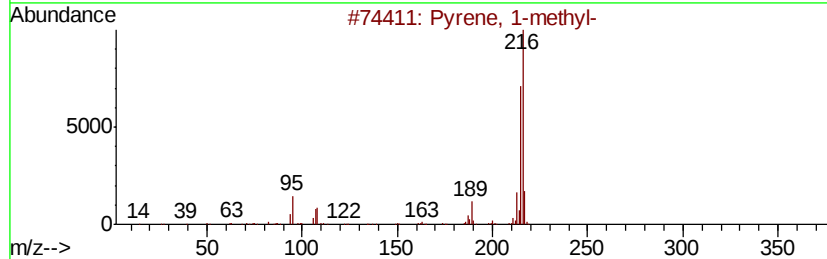
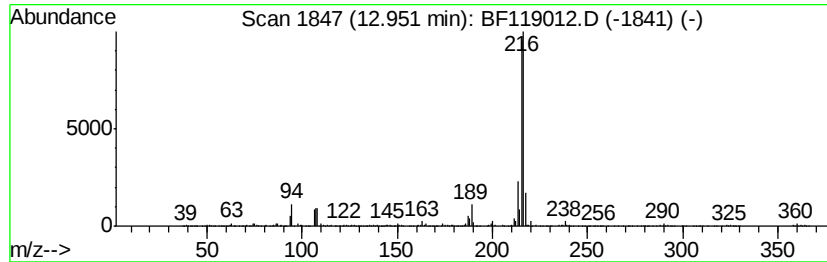
Instrument :  
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ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 15 Pyrene, 1-methyl- Concentration Rank 14

| R.T.    | EstConc | Area                    | Relative to ISTD | R.T.    |             |      |
|---------|---------|-------------------------|------------------|---------|-------------|------|
| 12.95   | 5.34 ng | 5106080                 | Chrysene-d12     | 13.95   |             |      |
| Hit# of | 5       | Tentative ID            | MW               | MolForm | CAS#        | Qual |
| 1       |         | Pyrene, 1-methyl-       | 216              | C17H12  | 002381-21-7 | 96   |
| 2       |         | Fluoranthene, 2-methyl- | 216              | C17H12  | 033543-31-6 | 91   |
| 3       |         | 11H-Benzo[b]fluorene    | 216              | C17H12  | 000243-17-4 | 90   |
| 4       |         | 11H-Benzo[a]fluorene    | 216              | C17H12  | 000238-84-6 | 83   |
| 5       |         | Pyrene, 2-methyl-       | 216              | C17H12  | 003442-78-2 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

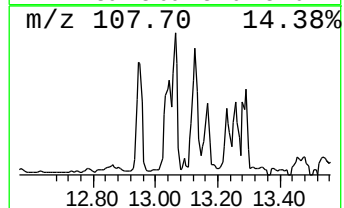
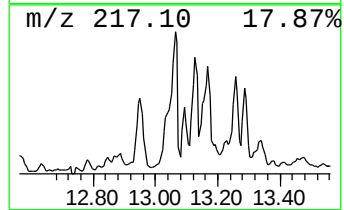
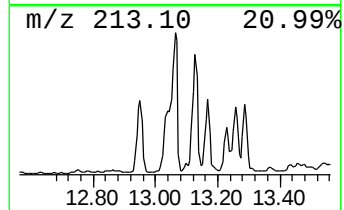
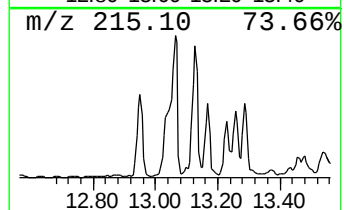
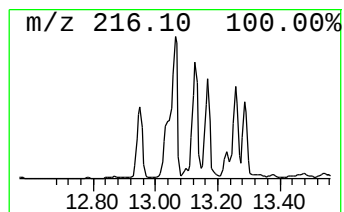
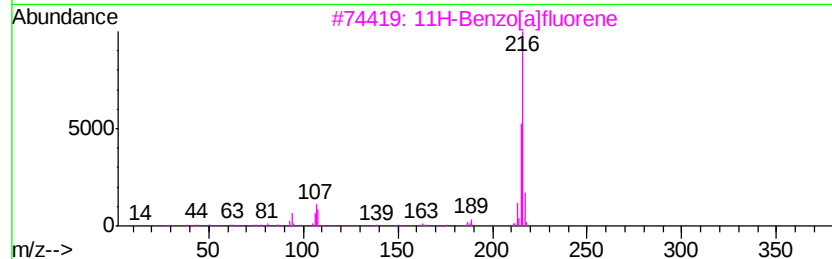
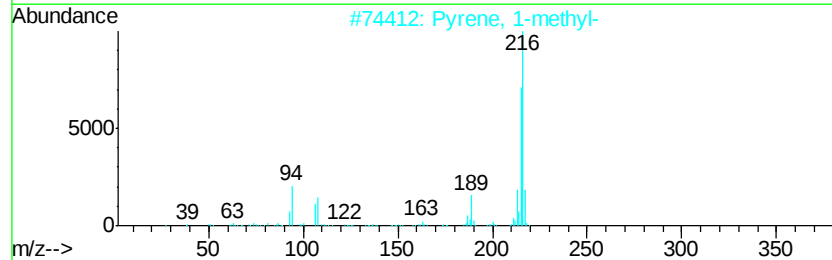
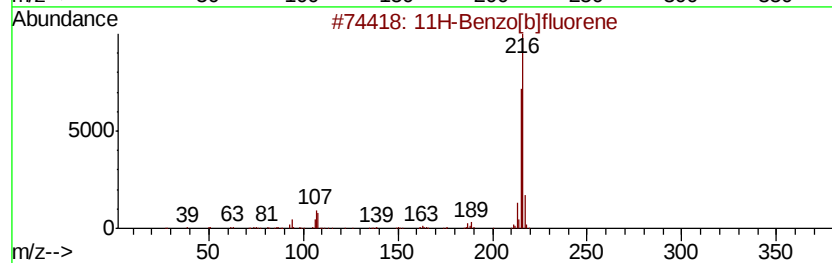
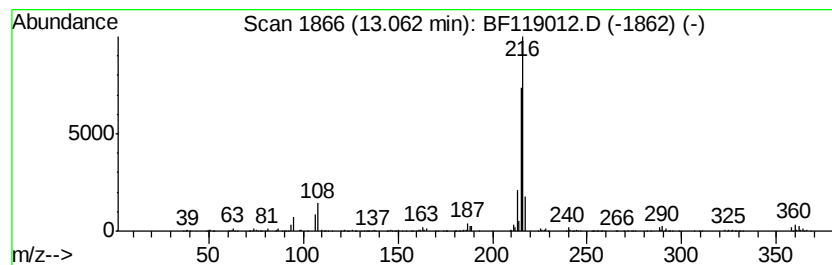
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 11H-Benzo[b]fluorene Concentration Rank 12

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 13.06 | 6.53 ng | 6239570 | Chrysene-d12     | 13.95 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 94   |
| 2    |    | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 91   |
| 3    |    | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 90   |
| 4    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 87   |
| 5    |    | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
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Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

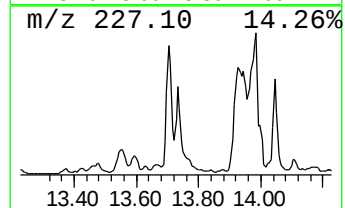
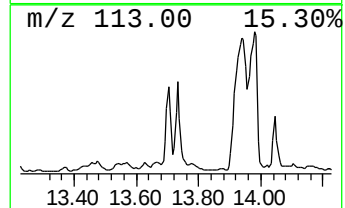
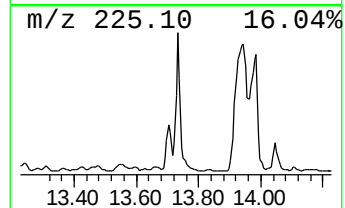
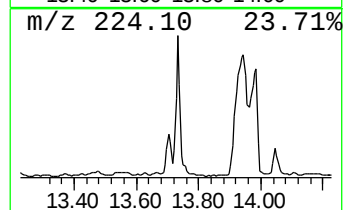
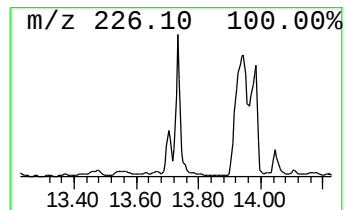
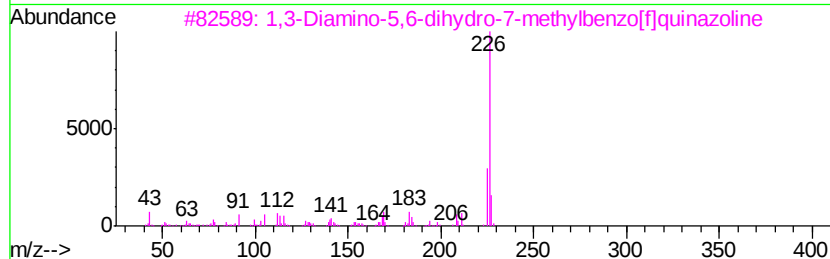
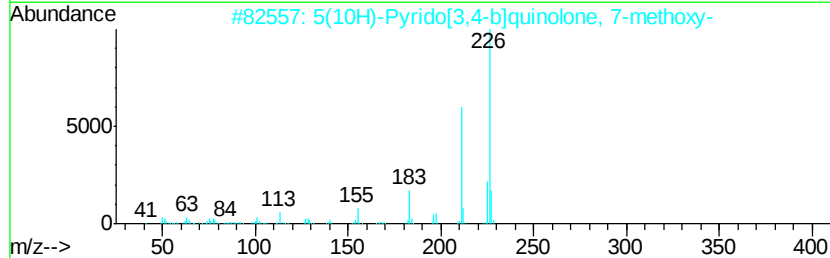
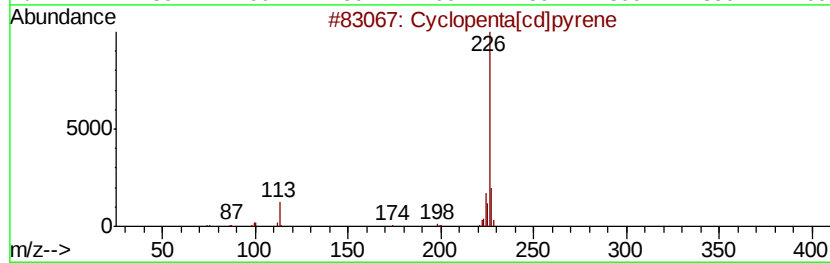
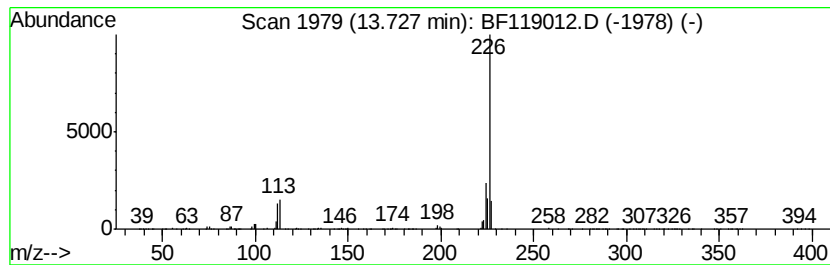
Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 17 Cyclopenta[cd]pyrene Concentration Rank 16

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.       |              |      |
|---------|---------|-------------------------------------|------------------|------------|--------------|------|
| 13.73   | 4.78 ng | 4567870                             | Chrysene-d12     | 13.95      |              |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm    | CAS#         | Qual |
| 1       |         | Cvclopenta[cd]pvrene                | 226              | C18H10     | 027208-37-3  | 76   |
| 2       |         | 5(10H)-Pyrido[3,4-b]quinolone, 7... | 226              | C13H10N2O2 | 1000256-99-5 | 53   |
| 3       |         | 1,3-Diamino-5,6-dihydro-7-methyl... | 226              | C13H14N4   | 037436-37-6  | 50   |
| 4       |         | Benzene, [4-(3-ethynylphenyl)-1,... | 226              | C18H10     | 1000115-87-3 | 50   |
| 5       |         | 2,2'-Oxydibenzaldehyde              | 226              | C14H10O3   | 049590-51-4  | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022120\  
Data File : BF119012.D  
Acq On : 22 Feb 2020 00:37  
Operator : CG/JU  
Sample : L1613-01 5X  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

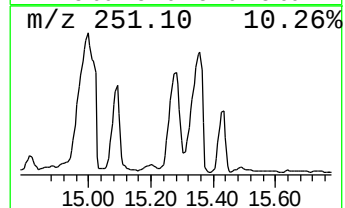
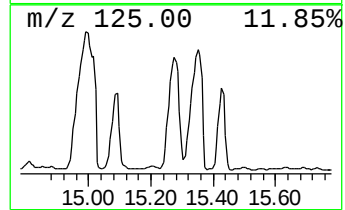
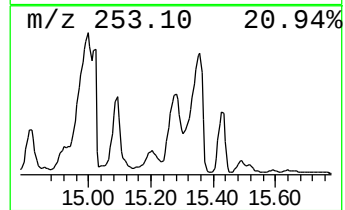
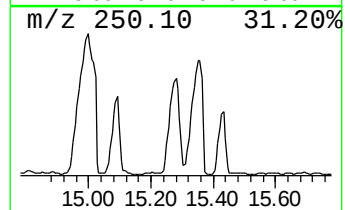
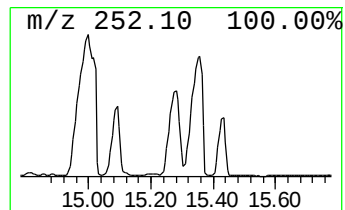
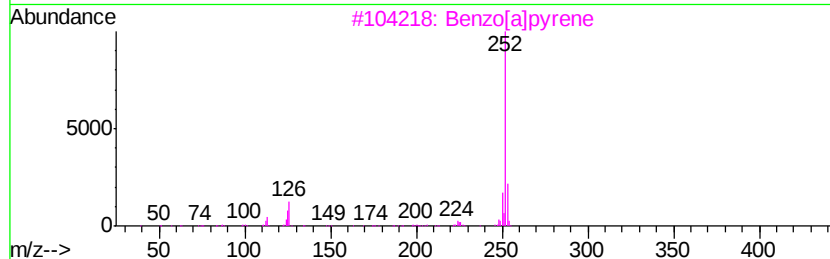
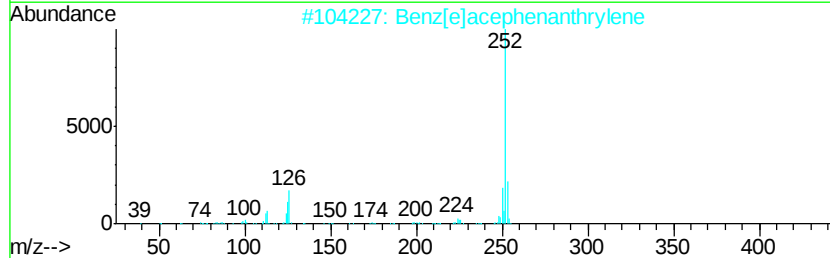
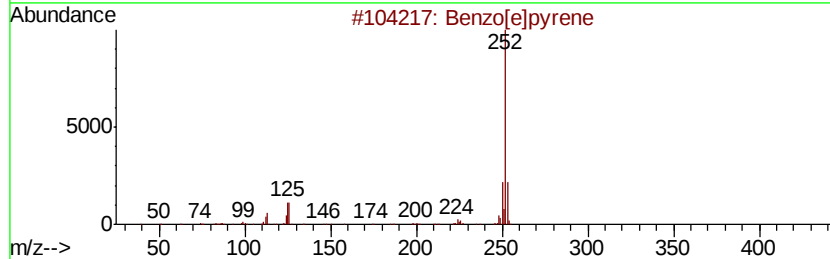
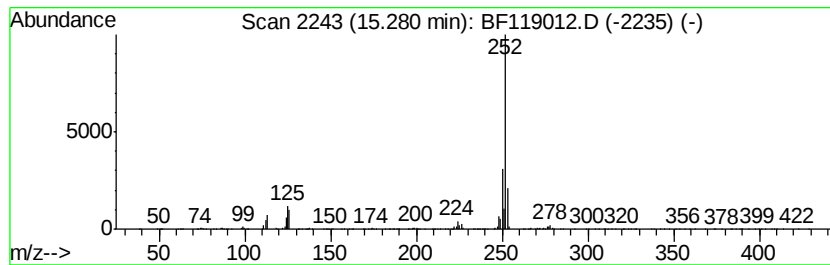
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Benzo[e]pyrene Concentration Rank 9

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 15.28 | 12.89 ng | 9194870 | Perylene-d12     | 15.39 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[e]bpvrene           | 252 | C20H12  | 000192-97-2 | 94   |
| 2    |    | Benzo[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 94   |
| 3    |    | Benzo[a]bpvrene           | 252 | C20H12  | 000050-32-8 | 93   |
| 4    |    | Benzo[i]lfluoranthene     | 252 | C20H12  | 000205-82-3 | 90   |
| 5    |    | Perylene                  | 252 | C20H12  | 000198-55-0 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF022120\  
 Data File : BF119012.D  
 Acq On : 22 Feb 2020 00:37  
 Operator : CG/JU  
 Sample : L1613-01 5X  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-14-EP1-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF022020.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 |
| Indene               | 7.02  | 15.8    | ng    | 634566   | 1                     | 6.76  | 801620   | 20.0 |
| Naphthalene, 1,6-... | 9.37  | 18.1    | ng    | 816490   | 3                     | 9.79  | 902685   | 20.0 |
| Naphthalene, 2,3-... | 9.45  | 21.4    | ng    | 968031   | 3                     | 9.79  | 902685   | 20.0 |
| Naphthalene, 2-et... | 9.52  | 21.6    | ng    | 974883   | 3                     | 9.79  | 902685   | 20.0 |
| Naphthalene, 2,3,... | 10.10 | 14.9    | ng    | 673340   | 3                     | 9.79  | 902685   | 20.0 |
| Dibenzofuran, 4-m... | 10.49 | 20.4    | ng    | 919181   | 3                     | 9.79  | 902685   | 20.0 |
| Anthracene, 1-met... | 11.78 | 4.0     | ng    | 2765360  | 4                     | 11.28 | 13931300 | 20.0 |
| Phenanthrene, 2-m... | 11.81 | 4.2     | ng    | 2922050  | 4                     | 11.28 | 13931300 | 20.0 |
| 4H-Cyclopenta[def... | 11.90 | 7.9     | ng    | 5529770  | 4                     | 11.28 | 13931300 | 20.0 |
| unknown12.37         | 12.37 | 4.1     | ng    | 2881160  | 4                     | 11.28 | 13931300 | 20.0 |
| Cyclopenta(def)ph... | 12.44 | 4.3     | ng    | 2995110  | 4                     | 11.28 | 13931300 | 20.0 |
| Pyrene, 1-methyl-    | 12.95 | 5.3     | ng    | 5106080  | 5                     | 13.95 | 19110200 | 20.0 |
| 11H-Benzo[b]fluorene | 13.06 | 6.5     | ng    | 6239570  | 5                     | 13.95 | 19110200 | 20.0 |
| Cyclopenta[cd]pyrene | 13.73 | 4.8     | ng    | 4567870  | 5                     | 13.95 | 19110200 | 20.0 |
| Benzo[e]pyrene       | 15.28 | 12.9    | ng    | 9194870  | 6                     | 15.39 | 14263100 | 20.0 |