

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
 Data File : BF111468.D  
 Acq On : 15 Dec 2018 15:03  
 Operator : JU/SJ  
 Sample : J6316-01  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BORING-SAMPLE-SIGNAL-TOWER-

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-BF121418.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 5.004       | 493           | 496         | 504          | rVB      | 164283         | 199310        | 4.78%           | 0.364%        |
| 2         | 5.422       | 562           | 567         | 570          | rBV      | 3943523        | 3637202       | 87.30%          | 6.634%        |
| 3         | 5.628       | 598           | 602         | 610          | rBV      | 238565         | 280337        | 6.73%           | 0.511%        |
| 4         | 6.351       | 722           | 725         | 728          | rVB      | 80179          | 64535         | 1.55%           | 0.118%        |
| 5         | 6.440       | 735           | 740         | 746          | rBV      | 3967361        | 4166125       | 100.00%         | 7.599%        |
| 6         | 6.569       | 757           | 762         | 765          | rBV      | 4445912        | 3894123       | 93.47%          | 7.103%        |
| 7         | 6.634       | 768           | 773         | 777          | rVV      | 298049         | 375770        | 9.02%           | 0.685%        |
| 8         | 6.669       | 777           | 779         | 785          | rVB      | 157148         | 164088        | 3.94%           | 0.299%        |
| 9         | 6.792       | 797           | 800         | 808          | rBV      | 1333771        | 1053372       | 25.28%          | 1.921%        |
| 10        | 6.951       | 823           | 827         | 830          | rBV      | 3681369        | 2997948       | 71.96%          | 5.468%        |
| 11        | 7.057       | 841           | 845         | 853          | rBV      | 441574         | 507757        | 12.19%          | 0.926%        |
| 12        | 7.351       | 889           | 895         | 901          | rBV      | 2405417        | 2582340       | 61.98%          | 4.710%        |
| 13        | 7.410       | 901           | 905         | 910          | rVV2     | 87152          | 142215        | 3.41%           | 0.259%        |
| 14        | 7.457       | 910           | 913         | 919          | rVB      | 77422          | 78828         | 1.89%           | 0.144%        |
| 15        | 7.569       | 926           | 932         | 934          | rBV3     | 39233          | 48943         | 1.17%           | 0.089%        |
| 16        | 7.651       | 940           | 946         | 956          | rBV      | 841135         | 845890        | 20.30%          | 1.543%        |
| 17        | 7.828       | 970           | 976         | 979          | rBV2     | 111202         | 133410        | 3.20%           | 0.243%        |
| 18        | 7.875       | 982           | 984         | 988          | rVB      | 148828         | 125941        | 3.02%           | 0.230%        |
| 19        | 8.075       | 1014          | 1018        | 1020         | rBV      | 1648991        | 1384944       | 33.24%          | 2.526%        |
| 20        | 8.098       | 1020          | 1022        | 1025         | rVB      | 457695         | 375391        | 9.01%           | 0.685%        |
| 21        | 8.292       | 1051          | 1055        | 1058         | rBV      | 747360         | 606983        | 14.57%          | 1.107%        |
| 22        | 8.351       | 1061          | 1065        | 1078         | rVV      | 789889         | 705849        | 16.94%          | 1.288%        |
| 23        | 8.781       | 1134          | 1138        | 1145         | rBV2     | 527703         | 579225        | 13.90%          | 1.057%        |
| 24        | 8.851       | 1146          | 1150        | 1153         | rVV      | 179028         | 160611        | 3.86%           | 0.293%        |
| 25        | 8.886       | 1153          | 1156        | 1159         | rVB2     | 186331         | 166626        | 4.00%           | 0.304%        |
| 26        | 8.963       | 1165          | 1169        | 1171         | rBV      | 132010         | 124493        | 2.99%           | 0.227%        |
| 27        | 9.069       | 1184          | 1187        | 1191         | rVB      | 202872         | 162085        | 3.89%           | 0.296%        |
| 28        | 9.122       | 1191          | 1196        | 1197         | rBV3     | 49175          | 47399         | 1.14%           | 0.086%        |
| 29        | 9.151       | 1197          | 1201        | 1204         | rBV      | 4780621        | 4158759       | 99.82%          | 7.586%        |
| 30        | 9.275       | 1220          | 1222        | 1225         | rBV      | 260512         | 226167        | 5.43%           | 0.413%        |
| 31        | 9.304       | 1225          | 1227        | 1232         | rVB      | 85813          | 78005         | 1.87%           | 0.142%        |
| 32        | 9.363       | 1232          | 1237        | 1239         | rVB2     | 39053          | 53226         | 1.28%           | 0.097%        |
| 33        | 9.392       | 1239          | 1242        | 1244         | rBV      | 44941          | 47427         | 1.14%           | 0.087%        |
| 34        | 9.486       | 1255          | 1258        | 1261         | rBV2     | 68590          | 71991         | 1.73%           | 0.131%        |

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

Filtering: 5  
 Min Area: 3 % of largest Peak  
 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-BF121418.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 9.539  | 1264 | 1267 | 1273 | rVB  | 352647  | 364437  | 8.75%  | 0.665% |
| 36 | 9.686  | 1288 | 1292 | 1298 | rBV  | 135367  | 143756  | 3.45%  | 0.262% |
| 37 | 9.769  | 1303 | 1306 | 1308 | rBV  | 123594  | 107033  | 2.57%  | 0.195% |
| 38 | 9.798  | 1308 | 1311 | 1313 | rVV  | 71710   | 98359   | 2.36%  | 0.179% |
| 39 | 9.828  | 1313 | 1316 | 1319 | rVV2 | 1627672 | 1447946 | 34.76% | 2.641% |
| 40 | 9.857  | 1319 | 1321 | 1330 | rVB  | 165627  | 192124  | 4.61%  | 0.350% |
| 41 | 10.033 | 1348 | 1351 | 1355 | rVB  | 64278   | 55277   | 1.33%  | 0.101% |
| 42 | 10.375 | 1406 | 1409 | 1416 | rVB  | 153357  | 146788  | 3.52%  | 0.268% |
| 43 | 10.439 | 1416 | 1420 | 1422 | rBV4 | 30903   | 46843   | 1.12%  | 0.085% |
| 44 | 10.616 | 1446 | 1450 | 1457 | rBV  | 2163109 | 1907650 | 45.79% | 3.480% |
| 45 | 10.739 | 1469 | 1471 | 1478 | rVB5 | 65091   | 71404   | 1.71%  | 0.130% |
| 46 | 10.922 | 1498 | 1502 | 1505 | rBV4 | 34763   | 51079   | 1.23%  | 0.093% |
| 47 | 11.116 | 1532 | 1535 | 1539 | rVB3 | 60977   | 75491   | 1.81%  | 0.138% |
| 48 | 11.169 | 1541 | 1544 | 1546 | rVV2 | 47715   | 60699   | 1.46%  | 0.111% |
| 49 | 11.210 | 1549 | 1551 | 1556 | rVB2 | 74845   | 79975   | 1.92%  | 0.146% |
| 50 | 11.310 | 1565 | 1568 | 1570 | rBV  | 1206121 | 1120570 | 26.90% | 2.044% |
| 51 | 11.333 | 1570 | 1572 | 1576 | rVV  | 1073543 | 946715  | 22.72% | 1.727% |
| 52 | 11.386 | 1578 | 1581 | 1585 | rVV  | 309581  | 262708  | 6.31%  | 0.479% |
| 53 | 11.422 | 1585 | 1587 | 1592 | rVV4 | 39082   | 49854   | 1.20%  | 0.091% |
| 54 | 11.610 | 1616 | 1619 | 1622 | rBV2 | 59595   | 49081   | 1.18%  | 0.090% |
| 55 | 11.816 | 1650 | 1654 | 1656 | rBV  | 180260  | 183435  | 4.40%  | 0.335% |
| 56 | 11.845 | 1656 | 1659 | 1664 | rVV2 | 862597  | 1011686 | 24.28% | 1.845% |
| 57 | 11.886 | 1664 | 1666 | 1669 | rVV  | 130520  | 140143  | 3.36%  | 0.256% |
| 58 | 11.927 | 1669 | 1673 | 1675 | rVV  | 291985  | 350086  | 8.40%  | 0.639% |
| 59 | 11.945 | 1675 | 1676 | 1680 | rVB  | 126570  | 103505  | 2.48%  | 0.189% |
| 60 | 12.110 | 1701 | 1704 | 1706 | rBV  | 148873  | 126695  | 3.04%  | 0.231% |
| 61 | 12.369 | 1745 | 1748 | 1751 | rBV  | 101374  | 102675  | 2.46%  | 0.187% |
| 62 | 12.404 | 1752 | 1754 | 1759 | rVB5 | 79954   | 100896  | 2.42%  | 0.184% |
| 63 | 12.451 | 1759 | 1762 | 1767 | rBV4 | 62319   | 80198   | 1.93%  | 0.146% |
| 64 | 12.527 | 1771 | 1775 | 1778 | rBV  | 1332071 | 1133557 | 27.21% | 2.068% |
| 65 | 12.633 | 1788 | 1793 | 1798 | rBV3 | 612006  | 737116  | 17.69% | 1.345% |
| 66 | 12.751 | 1808 | 1813 | 1816 | rBV  | 1295284 | 1428300 | 34.28% | 2.605% |
| 67 | 12.804 | 1820 | 1822 | 1827 | rVB2 | 99885   | 105624  | 2.54%  | 0.193% |
| 68 | 12.898 | 1834 | 1838 | 1841 | rBV  | 2939891 | 2473197 | 59.36% | 4.511% |
| 69 | 12.974 | 1848 | 1851 | 1854 | rVB2 | 81418   | 100094  | 2.40%  | 0.183% |
| 70 | 13.057 | 1863 | 1865 | 1866 | rBV2 | 61118   | 54852   | 1.32%  | 0.100% |
| 71 | 13.080 | 1867 | 1869 | 1872 | rVB2 | 217454  | 179783  | 4.32%  | 0.328% |

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 BORING-SAMPLE-SIGNAL-TOWER-

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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF121418.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 72 | 13.145 | 1878 | 1880 | 1883 | rVB  | 182551  | 161667  | 3.88%  | 0.295% |
| 73 | 13.186 | 1884 | 1887 | 1889 | rBV  | 148323  | 139192  | 3.34%  | 0.254% |
| 74 | 13.274 | 1900 | 1902 | 1905 | rBV  | 94614   | 81880   | 1.97%  | 0.149% |
| 75 | 13.310 | 1905 | 1908 | 1912 | rVB  | 112840  | 101117  | 2.43%  | 0.184% |
| 76 | 13.727 | 1977 | 1979 | 1981 | rBV  | 121806  | 86369   | 2.07%  | 0.158% |
| 77 | 13.751 | 1981 | 1983 | 1987 | rVV2 | 116025  | 107261  | 2.57%  | 0.196% |
| 78 | 13.804 | 1990 | 1992 | 1998 | rVB2 | 157473  | 219336  | 5.26%  | 0.400% |
| 79 | 13.951 | 2012 | 2017 | 2019 | rBV2 | 1276613 | 1777311 | 42.66% | 3.242% |
| 80 | 13.974 | 2019 | 2021 | 2027 | rVB  | 788717  | 733440  | 17.60% | 1.338% |
| 81 | 14.057 | 2032 | 2035 | 2039 | rVB  | 306624  | 244932  | 5.88%  | 0.447% |
| 82 | 14.339 | 2080 | 2083 | 2086 | rVB  | 152020  | 142791  | 3.43%  | 0.260% |
| 83 | 14.368 | 2086 | 2088 | 2092 | rVB2 | 96149   | 88678   | 2.13%  | 0.162% |
| 84 | 14.427 | 2096 | 2098 | 2101 | rVV3 | 147001  | 148142  | 3.56%  | 0.270% |
| 85 | 14.463 | 2101 | 2104 | 2105 | rVV  | 108547  | 107097  | 2.57%  | 0.195% |
| 86 | 14.486 | 2105 | 2108 | 2111 | rVB  | 157347  | 162962  | 3.91%  | 0.297% |
| 87 | 14.980 | 2187 | 2192 | 2201 | rBV  | 664444  | 1334844 | 32.04% | 2.435% |
| 88 | 15.080 | 2207 | 2209 | 2213 | rVB  | 171677  | 153651  | 3.69%  | 0.280% |
| 89 | 15.268 | 2238 | 2241 | 2247 | rVB  | 420177  | 541342  | 12.99% | 0.987% |
| 90 | 15.333 | 2247 | 2252 | 2258 | rBV  | 722109  | 872706  | 20.95% | 1.592% |
| 91 | 15.392 | 2258 | 2262 | 2265 | rBV2 | 679834  | 855333  | 20.53% | 1.560% |
| 92 | 16.809 | 2498 | 2503 | 2514 | rVB2 | 239929  | 451067  | 10.83% | 0.823% |
| 93 | 17.233 | 2570 | 2575 | 2586 | rVB  | 171852  | 374738  | 8.99%  | 0.684% |

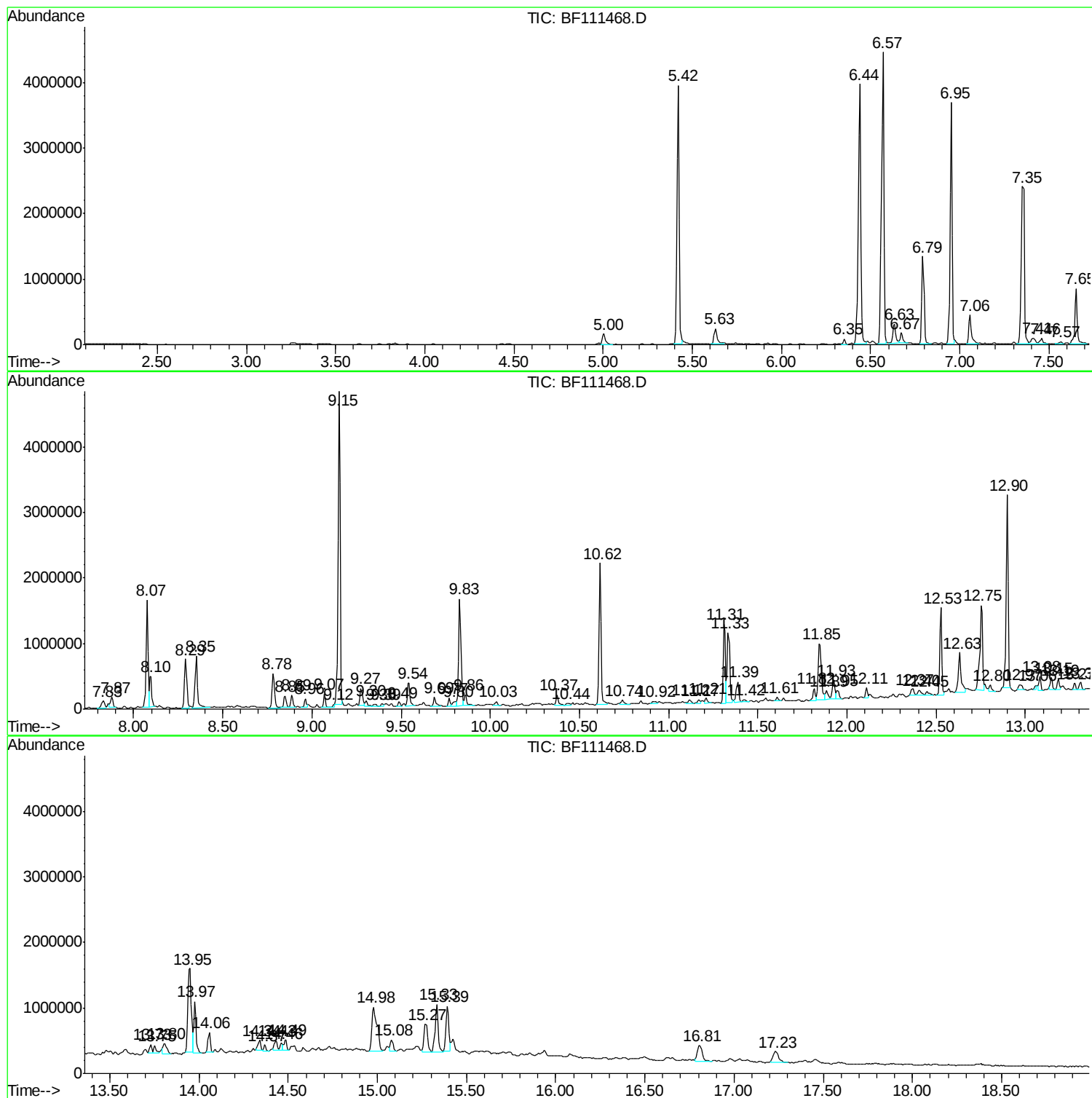
Sum of corrected areas: 54822802

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.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\BF121518\  
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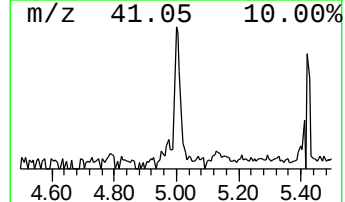
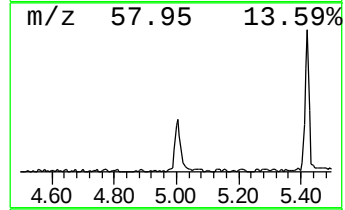
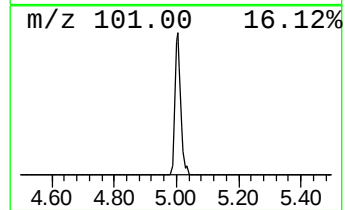
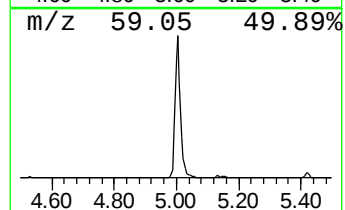
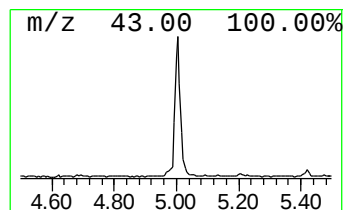
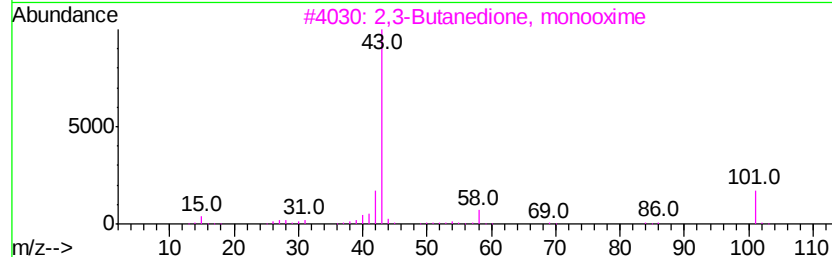
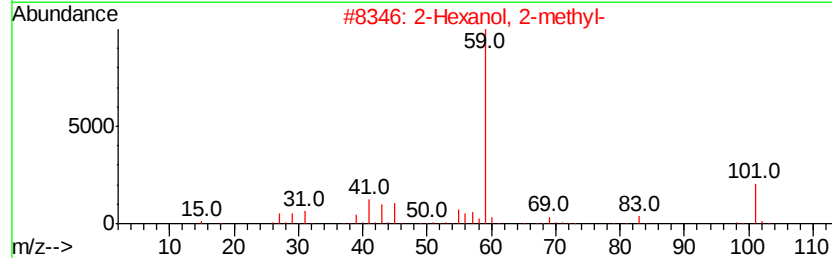
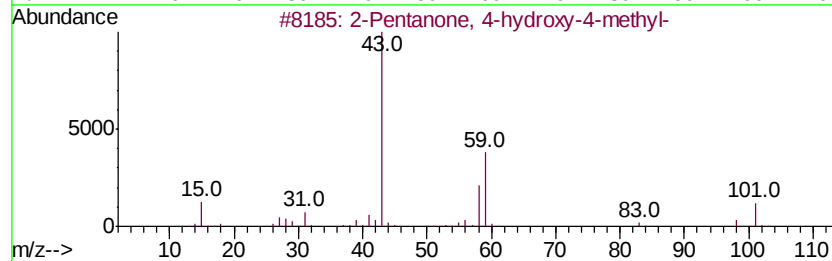
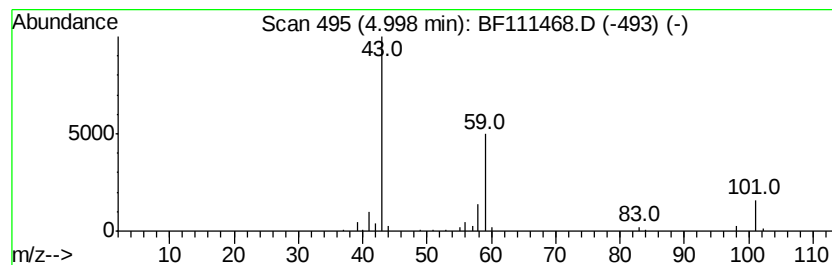
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.00 | 3.78 ng | 199310 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2 | 000123-42-2  | 64   |
| 2    |    |   | 2-Hexanol, 2-methyl-                | 116 | C7H16O  | 000625-23-0  | 28   |
| 3    |    |   | 2,3-Butanedione, monooxime          | 101 | C4H7NO2 | 000057-71-6  | 9    |
| 4    |    |   | n-Propyl t-butyl ether              | 116 | C7H16O  | 1000334-81-9 | 9    |
| 5    |    |   | Propane, 2-methyl-2-(1-methyleth... | 116 | C7H16O  | 017348-59-3  | 9    |



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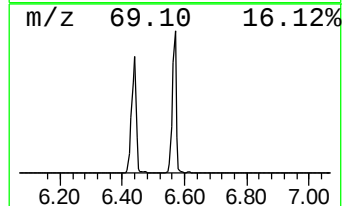
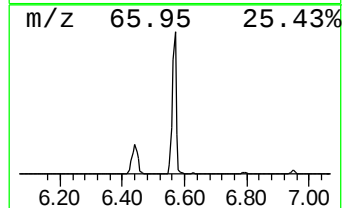
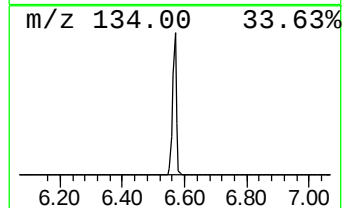
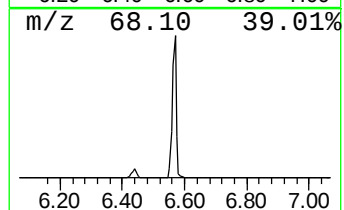
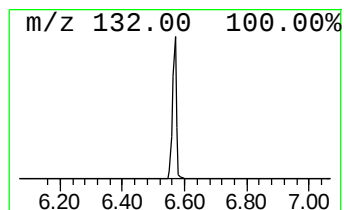
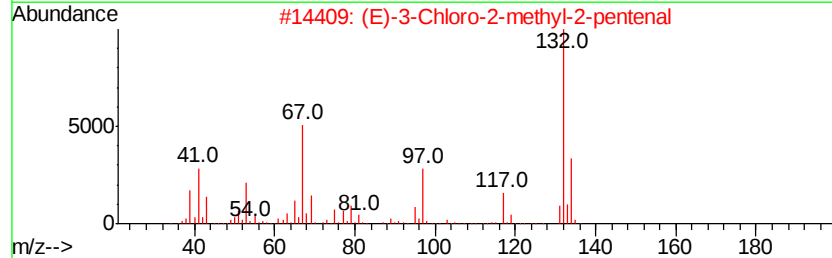
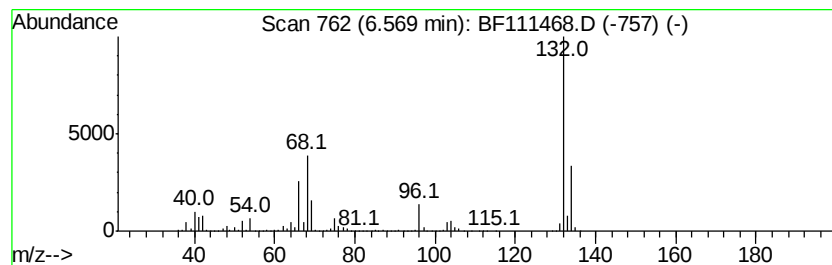
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 unknown6.57 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.57 | 73.94 ng | 3894120 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 16   |
| 3    |    | Tranvlcypropine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 4    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2    | 022291-85-6  | 9    |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

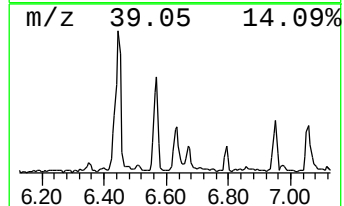
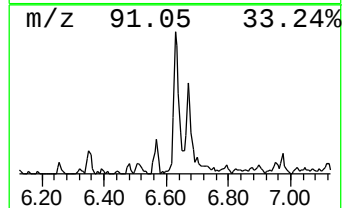
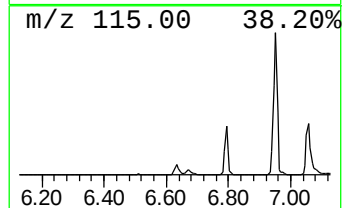
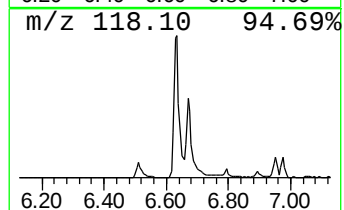
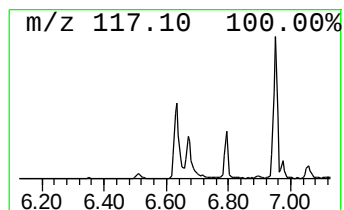
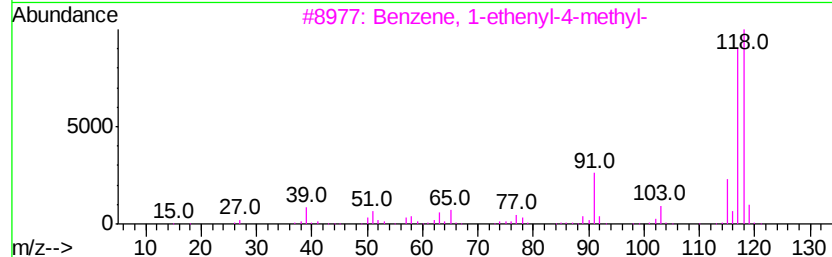
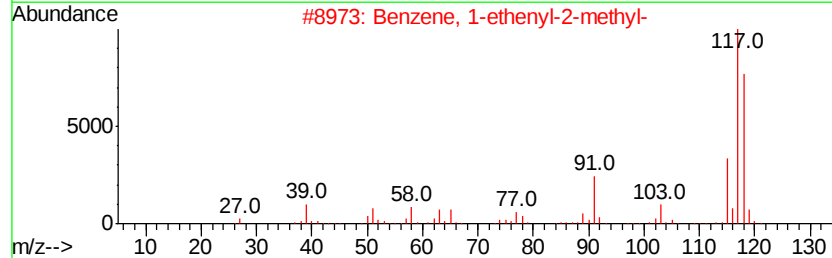
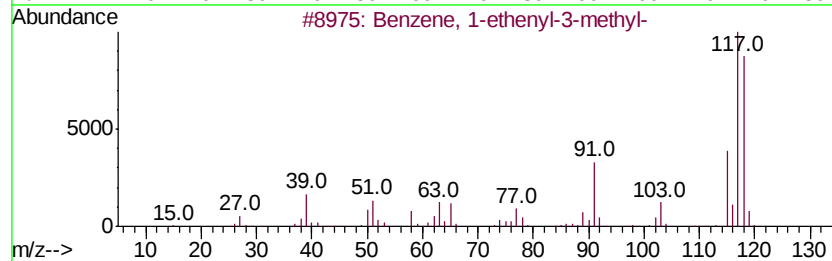
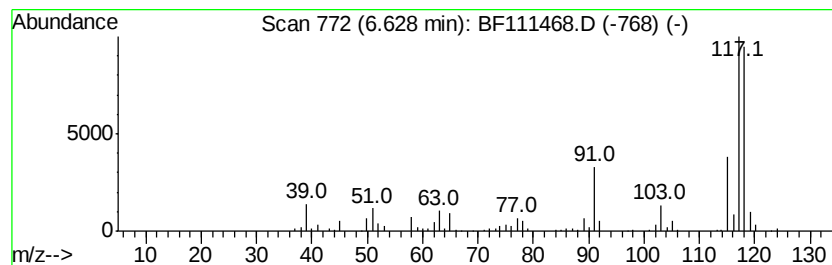
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Benzene, 1-ethenyl-3-methyl- Concentration Rank 10

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.63 | 7.13 ng | 375770 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 96   |
| 2    |    | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10   | 000611-15-4 | 94   |
| 3    |    | Benzene, 1-ethenyl-4-methyl- | 118 | C9H10   | 000622-97-9 | 93   |
| 4    |    | (Z)-1-Phenylpropene          | 118 | C9H10   | 000766-90-5 | 90   |
| 5    |    | Benzene, 2-propenyl-         | 118 | C9H10   | 000300-57-2 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

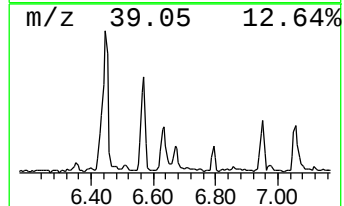
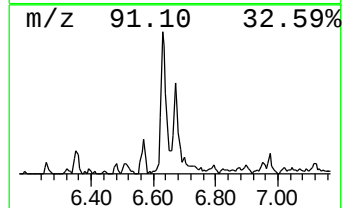
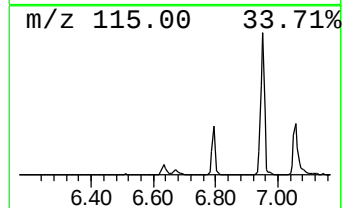
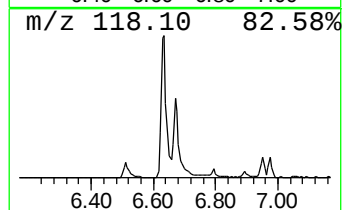
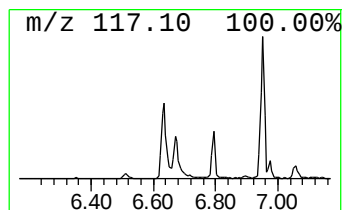
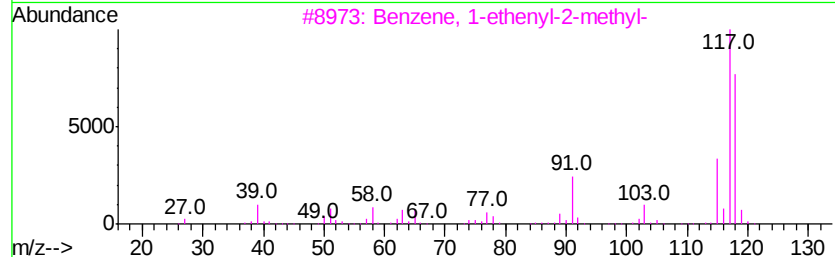
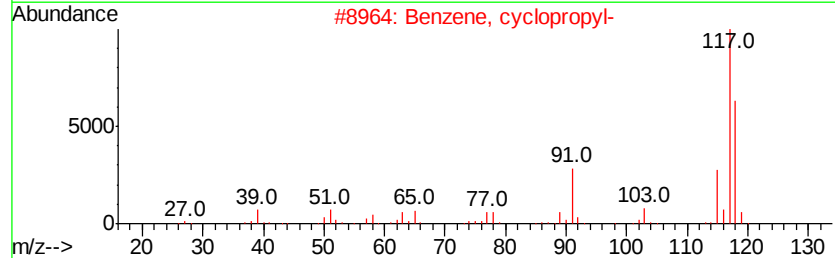
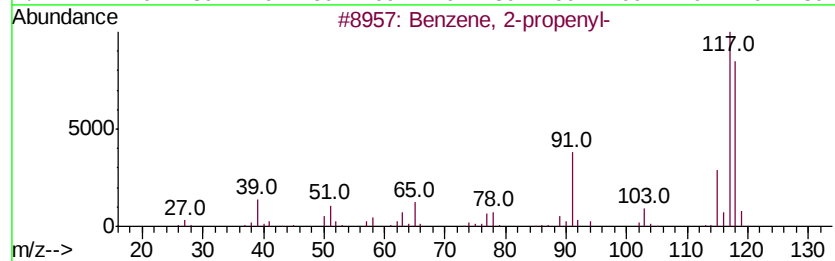
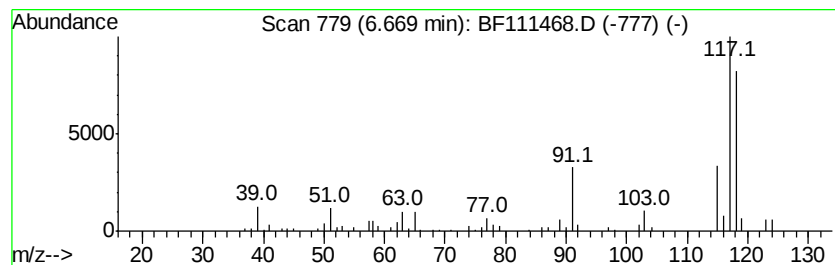
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 Benzene, 2-propenyl- Concentration Rank 17

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.67 | 3.12 ng | 164088 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-propenyl-                | 118 | C9H10   | 000300-57-2 | 94   |
| 2    |    | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4 | 94   |
| 3    |    | Benzene, 1-ethenyl-2-methyl-        | 118 | C9H10   | 000611-15-4 | 93   |
| 4    |    | Benzene, 1,1'-(1-ethenyl-1,3-pro... | 222 | C17H18  | 061141-97-7 | 91   |
| 5    |    | Benzene, 1-propenyl-                | 118 | C9H10   | 000637-50-3 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

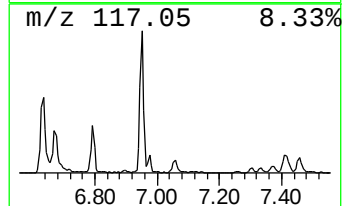
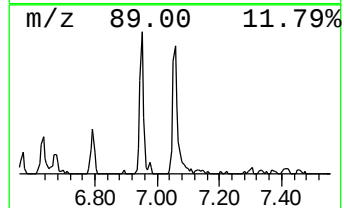
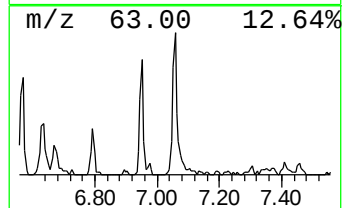
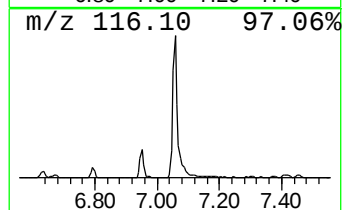
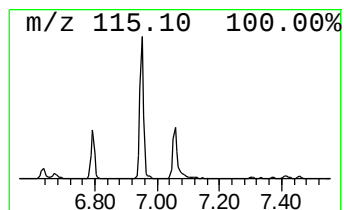
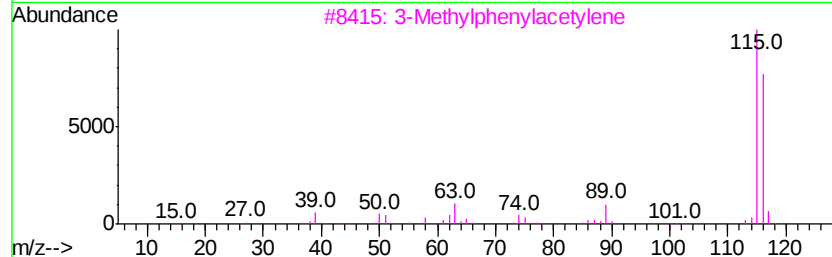
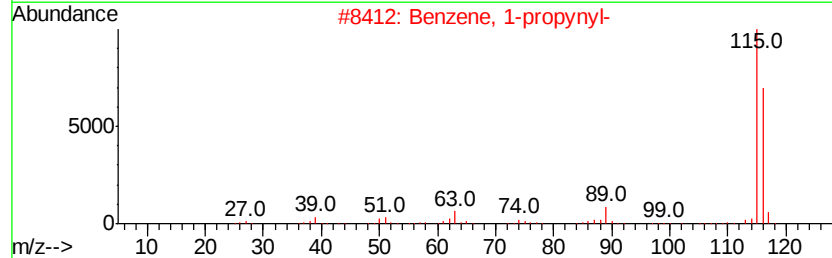
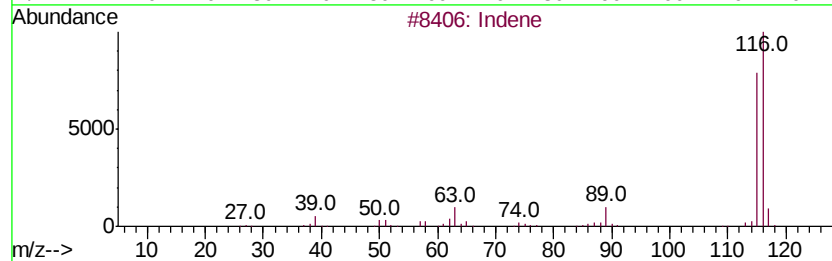
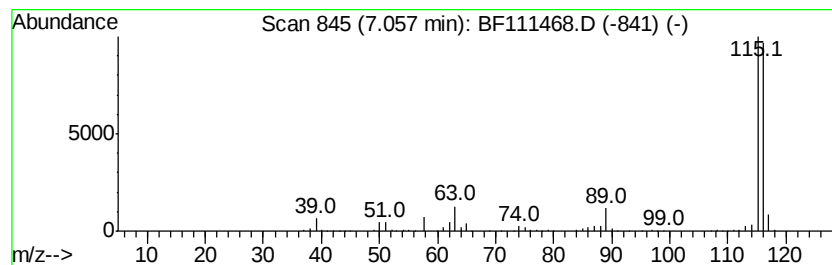
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Indene Concentration Rank 7

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.06 | 9.64 ng | 507757 | 1,4-Dichlorobenzene-d4 | 6.79 |

| 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 | 94   |
| 3    |    | 3-Methylphenylacetylene      | 116 | C9H8    | 000766-82-5 | 91   |
| 4    |    | Benzene, 1-ethynyl-4-methyl- | 116 | C9H8    | 000766-97-2 | 91   |
| 5    |    | 2-Methylphenylacetylene      | 116 | C9H8    | 000766-47-2 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

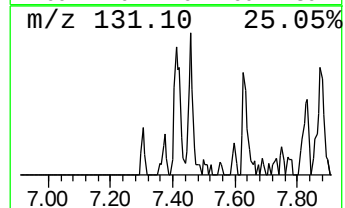
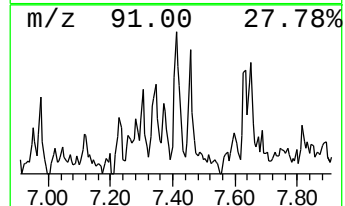
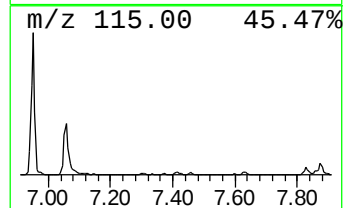
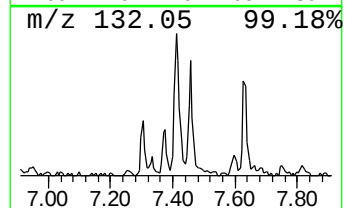
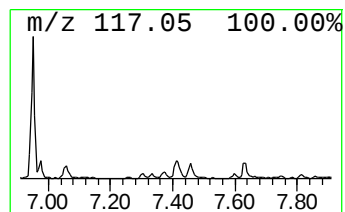
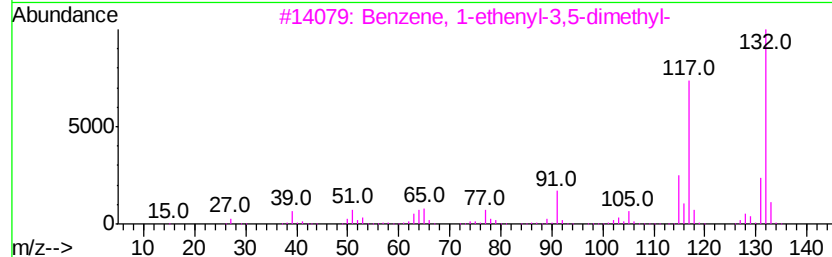
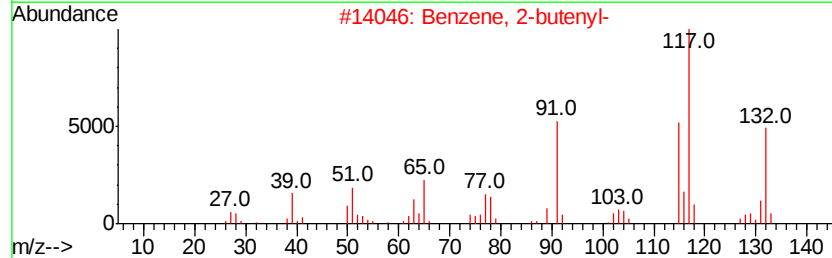
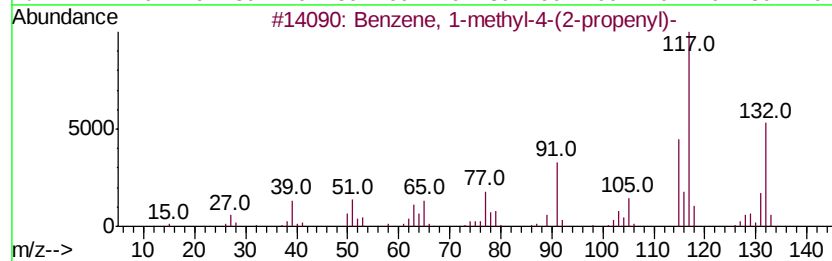
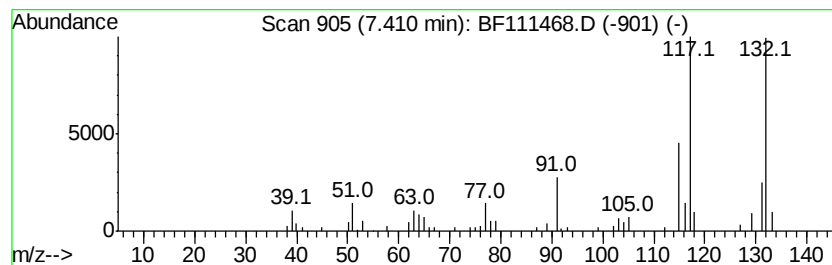
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Benzene, 1-methyl-4-(2-prop... Concentration Rank 19

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.41 | 2.70 ng | 142215 | 1,4-Dichlorobenzene-d4 | 6.79 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-4-(2-propenyl)- | 132 | C10H12  | 003333-13-9 | 97   |
| 2    |    | Benzene, 2-butenyl-               | 132 | C10H12  | 001560-06-1 | 95   |
| 3    |    | Benzene, 1-ethenyl-3,5-dimethyl-  | 132 | C10H12  | 005379-20-4 | 94   |
| 4    |    | Benzene, (2-methyl-1-propenyl)-   | 132 | C10H12  | 000768-49-0 | 94   |
| 5    |    | (E)-1-Phenyl-1-butene             | 132 | C10H12  | 001005-64-7 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

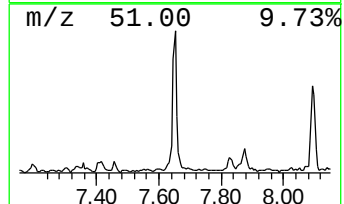
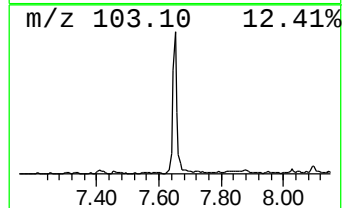
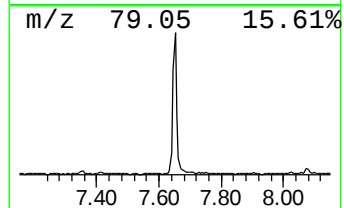
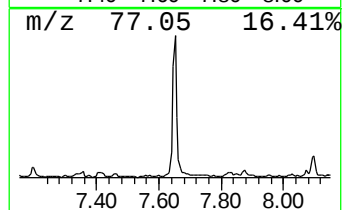
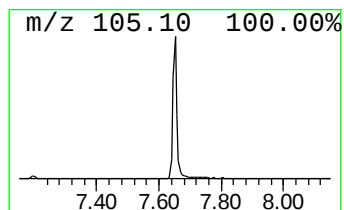
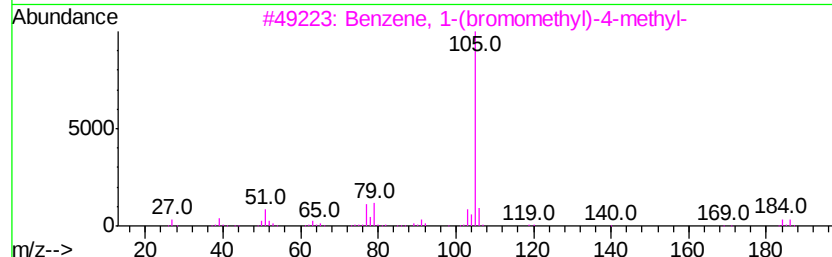
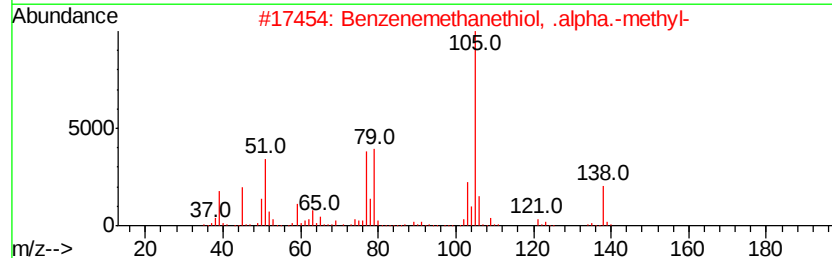
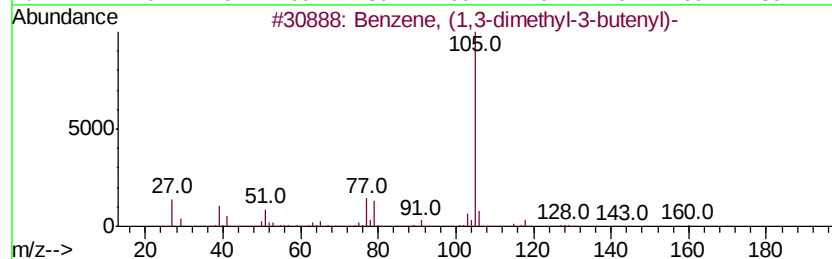
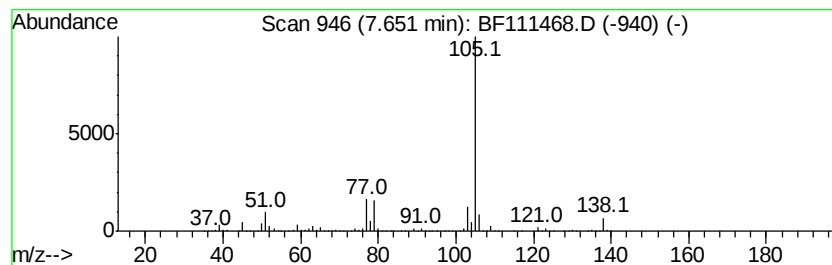
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Benzene, (1,3-dimethyl-3-bu... Concentration Rank 5

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 7.65 | 12.22 ng | 845890 | Naphthalene-d8   | 8.07 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------------------|-----|---------|-------------|------|
| 1       |   | Benzene, (1,3-dimethyl-3-butenyl)-  | 160 | C12H16  | 056851-51-5 | 72   |
| 2       |   | Benzenemethanethiol, .alpha.-met... | 138 | C8H10S  | 006263-65-6 | 72   |
| 3       |   | Benzene, 1-(bromomethyl)-4-methyl-  | 184 | C8H9Br  | 000104-81-4 | 72   |
| 4       |   | Benzenemethanethiol, 4-methyl-      | 138 | C8H10S  | 004498-99-1 | 64   |
| 5       |   | Ether, p-methylbenzyl vinyl         | 148 | C10H12O | 026437-10-5 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

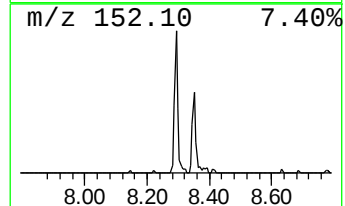
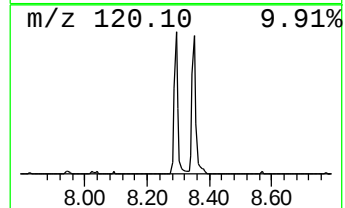
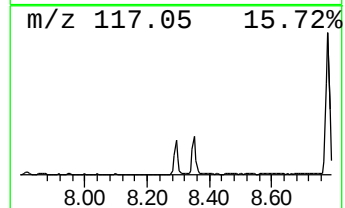
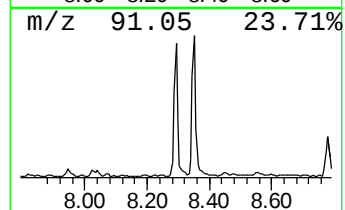
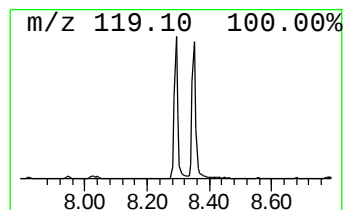
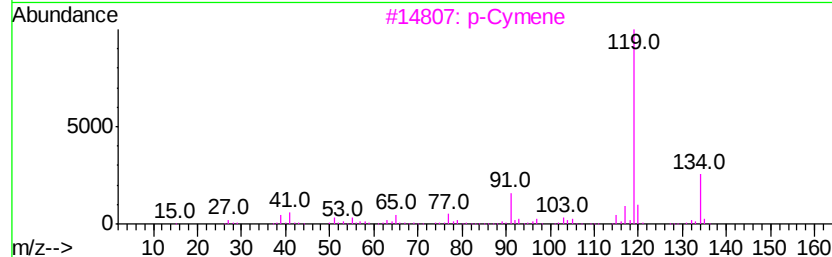
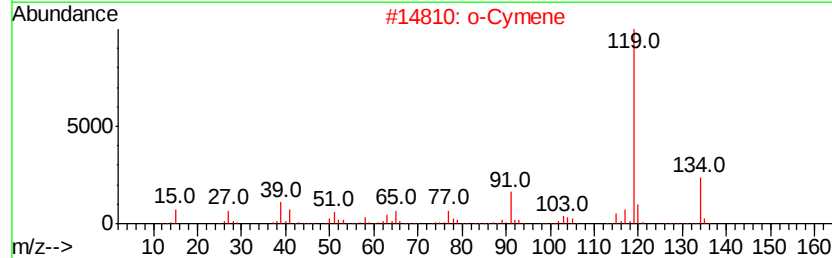
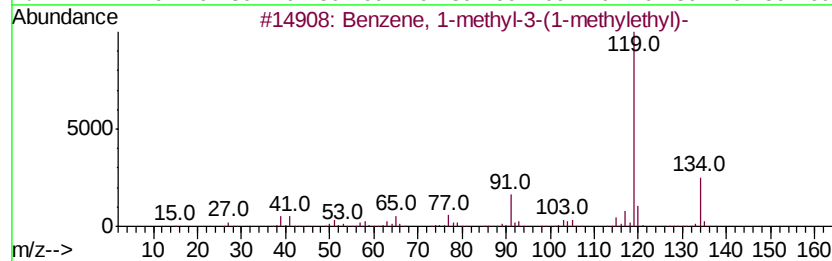
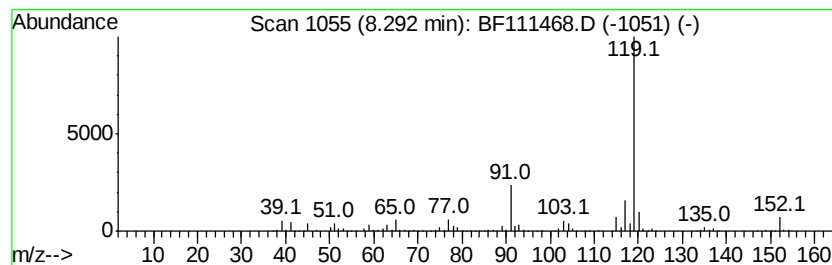
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Benzene, 1-methyl-3-(1-meth... Concentration Rank 8

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 8.29 | 8.77 ng | 606983 | Napthalene-d8    | 8.07 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 87   |
| 2    |    | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 81   |
| 3    |    | p-Cymene                            | 134 | C10H14  | 000099-87-6 | 80   |
| 4    |    | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0 | 72   |
| 5    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

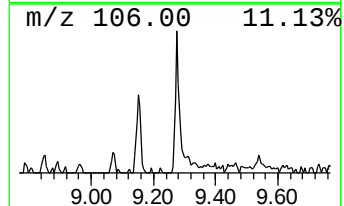
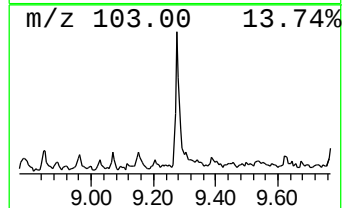
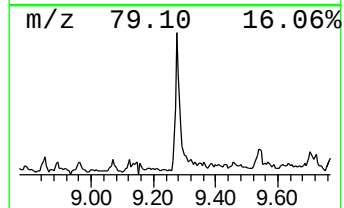
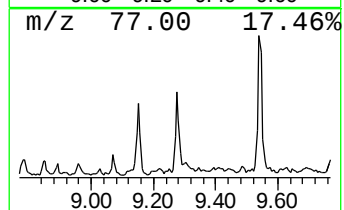
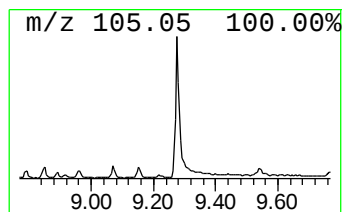
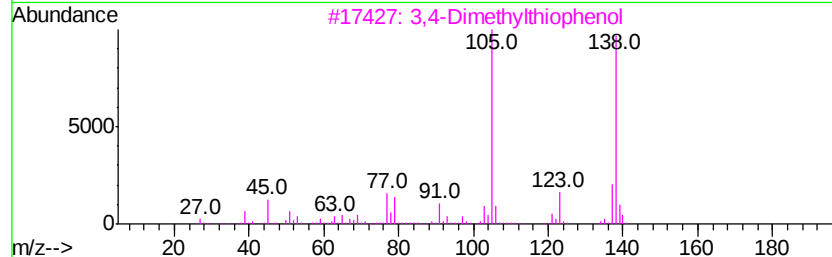
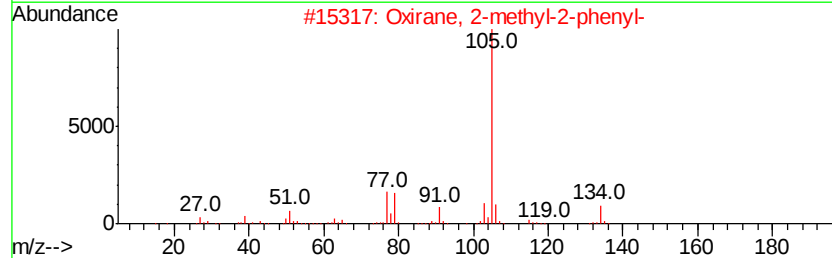
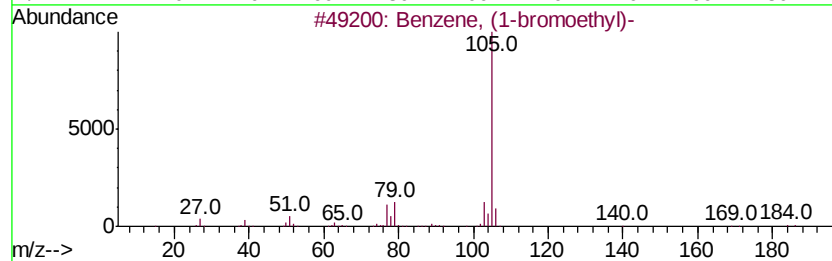
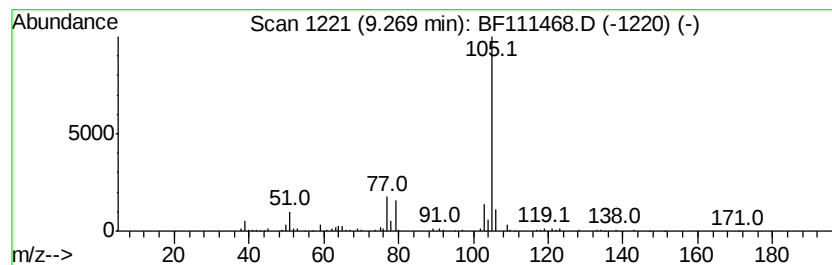
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Benzene, (1-bromoethyl)- Concentration Rank 16

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.27 | 3.12 ng | 226167 | Acenaphthene-d10 | 9.83 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (1-bromoethyl)-    | 184 | C8H9Br  | 000585-71-7 | 80   |
| 2    |    | Oxirane, 2-methyl-2-phenyl- | 134 | C9H10O  | 002085-88-3 | 72   |
| 3    |    | 3,4-Dimethylthiophenol      | 138 | C8H10S  | 018800-53-8 | 64   |
| 4    |    | Ether, p-methylbenzyl vinyl | 148 | C10H12O | 026437-10-5 | 64   |
| 5    |    | 3-Methylbenzyl mercaptan    | 138 | C8H10S  | 025697-56-7 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

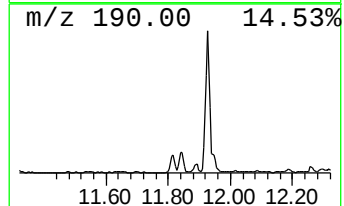
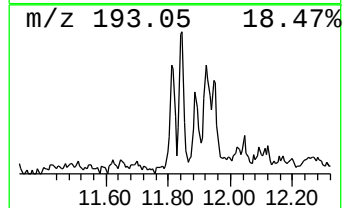
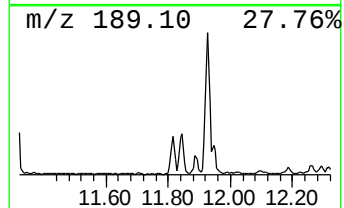
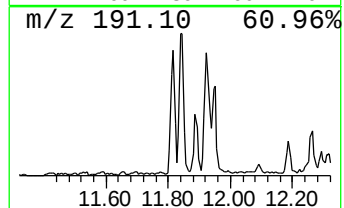
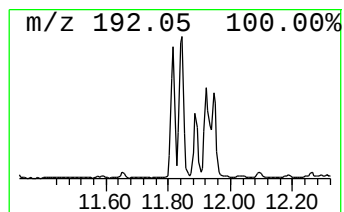
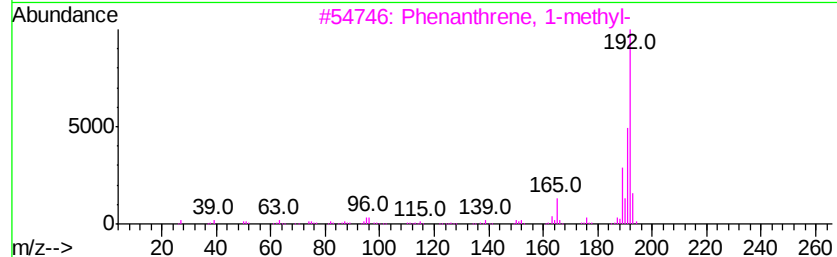
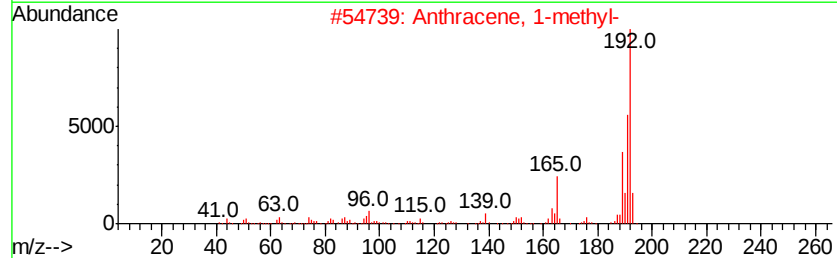
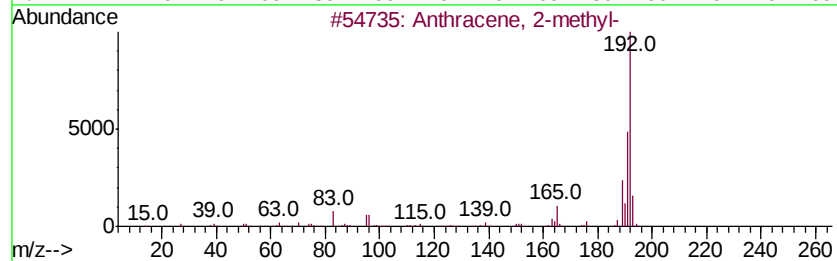
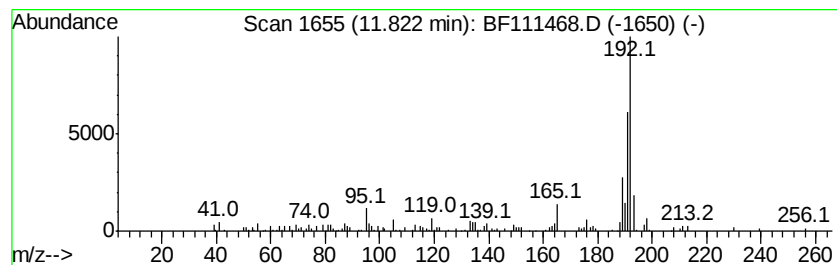
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Anthracene, 2-methyl- Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.82 | 3.27 ng | 183435 | Phenanthrene-d10 | 11.31 |

| Hit# | of | 5 | Tentative ID                           | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 2-methyl-                  | 192 | C15H12  | 000613-12-7 | 94   |
| 2    |    |   | Anthracene, 1-methyl-                  | 192 | C15H12  | 000610-48-0 | 94   |
| 3    |    |   | Phenanthrene, 1-methyl-                | 192 | C15H12  | 000832-69-9 | 93   |
| 4    |    |   | 1H-Cyclopropa[1,1]phenanthrene, 1a,... | 192 | C15H12  | 000949-41-7 | 90   |
| 5    |    |   | Phenanthrene, 2-methyl-                | 192 | C15H12  | 002531-84-2 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

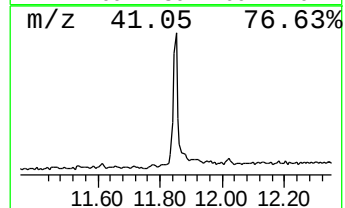
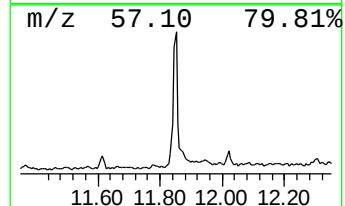
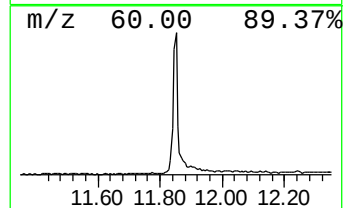
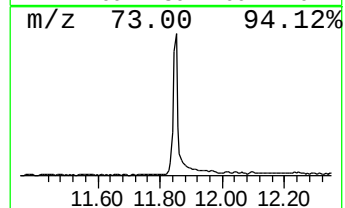
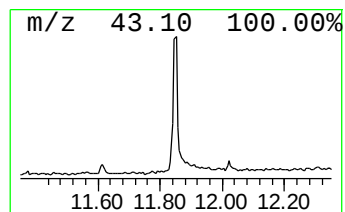
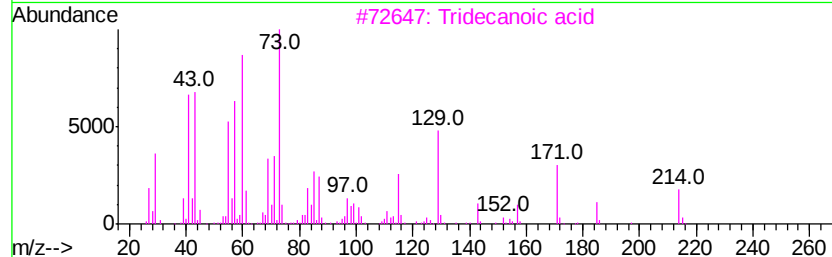
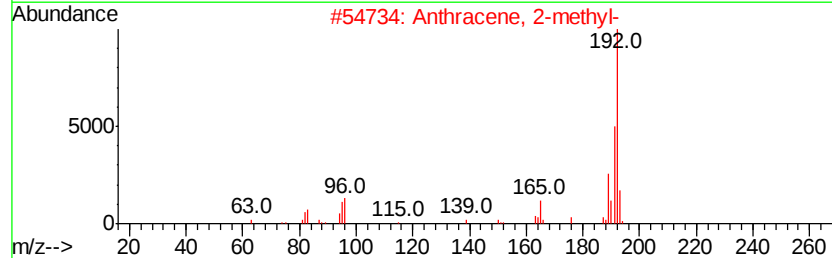
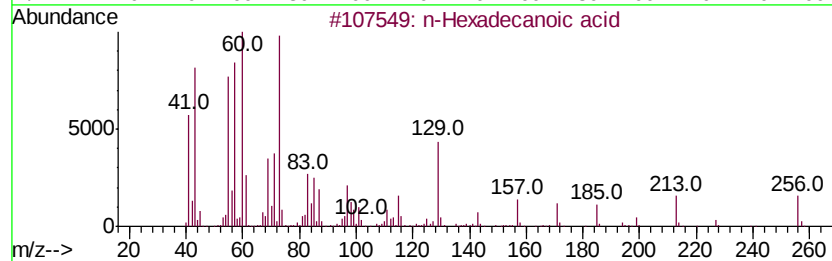
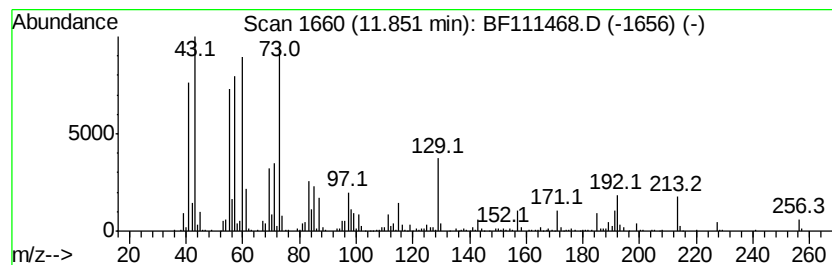
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 n-Hexadecanoic acid Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 11.85 | 18.06 ng | 1011690 | Phenanthrene-d10 | 11.31 |

| Hit# | of | Tentative ID                           | MW  | MolForm  | CAS#        | Qual |
|------|----|----------------------------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid                    | 256 | C16H32O2 | 000057-10-3 | 97   |
| 2    |    | Anthracene, 2-methyl-                  | 192 | C15H12   | 000613-12-7 | 86   |
| 3    |    | Tridecanoic acid                       | 214 | C13H26O2 | 000638-53-9 | 83   |
| 4    |    | 1H-Cyclopropa[1,1]phenanthrene, 1a,... | 192 | C15H12   | 000949-41-7 | 59   |
| 5    |    | Phenanthrene, 2-methyl-                | 192 | C15H12   | 002531-84-2 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

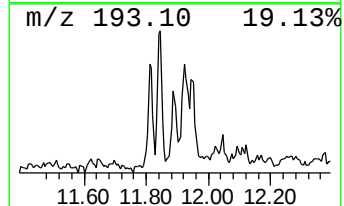
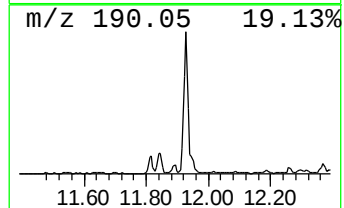
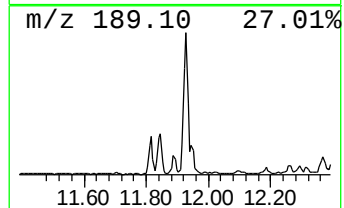
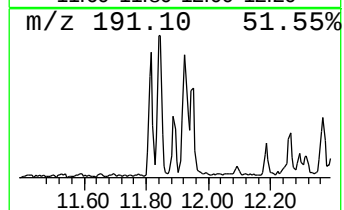
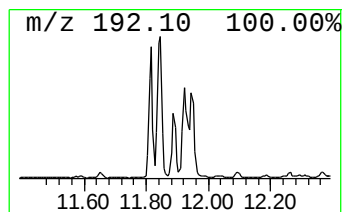
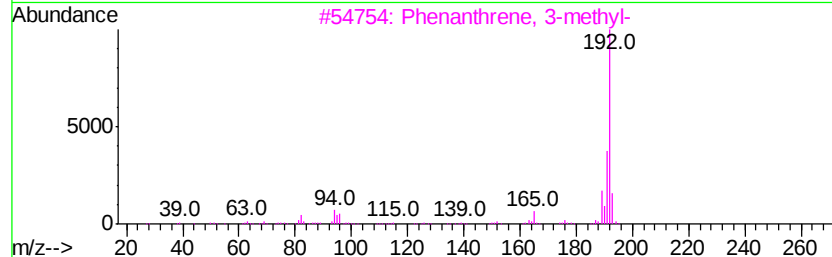
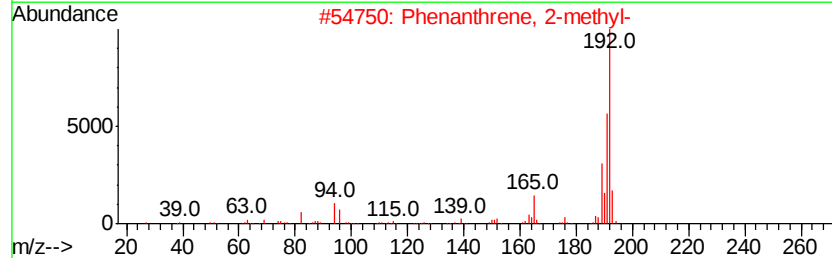
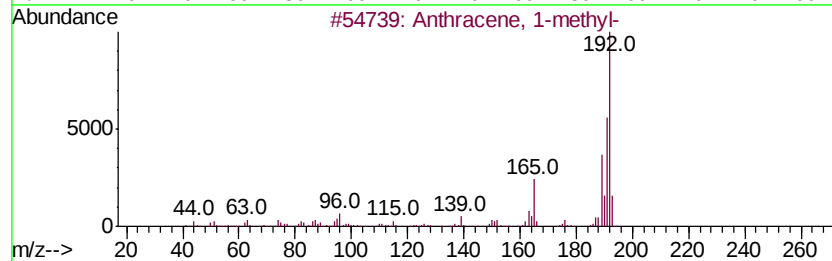
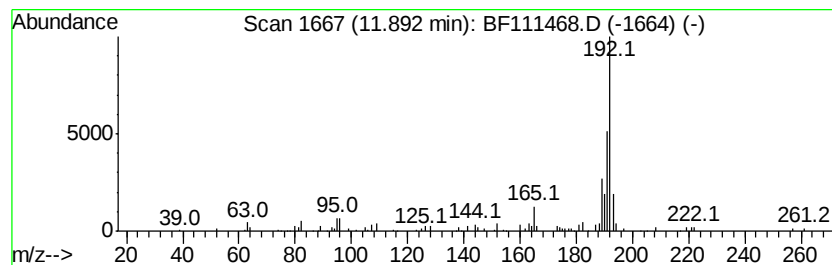
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Anthracene, 1-methyl- Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.89 | 2.50 ng | 140143 | Phenanthrene-d10 | 11.31 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 93   |
| 2    |    | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 91   |
| 3    |    | Phenanthrene, 3-methyl- | 192 | C15H12  | 000832-71-3 | 87   |
| 4    |    | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 87   |
| 5    |    | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

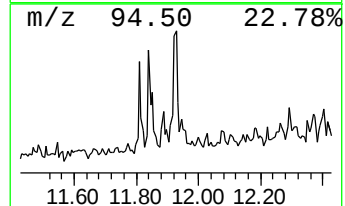
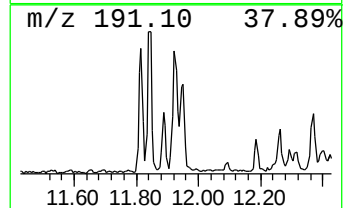
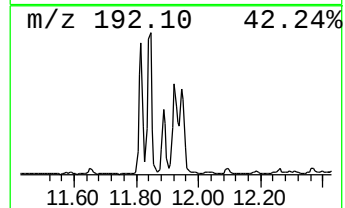
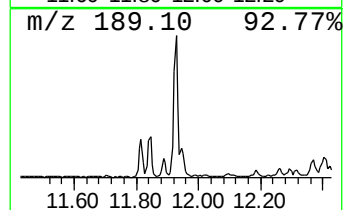
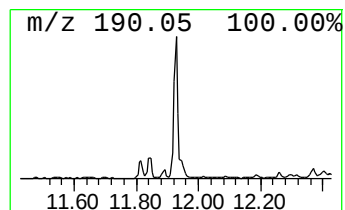
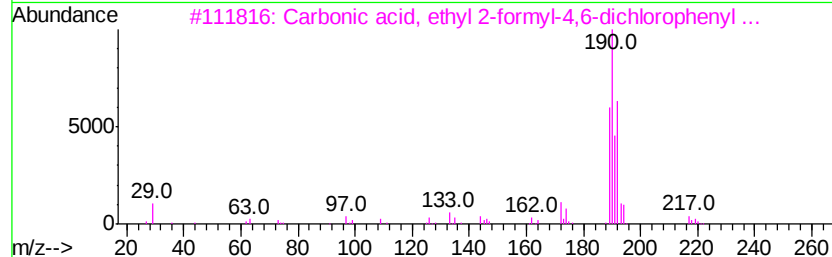
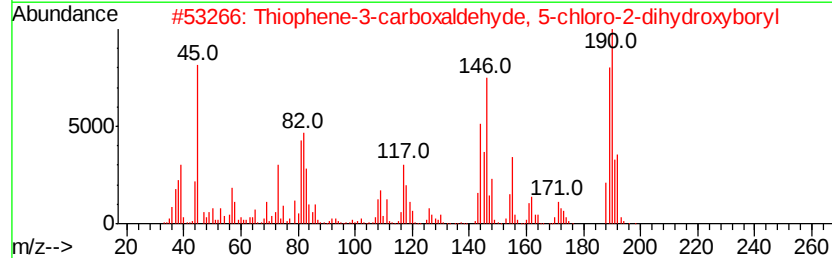
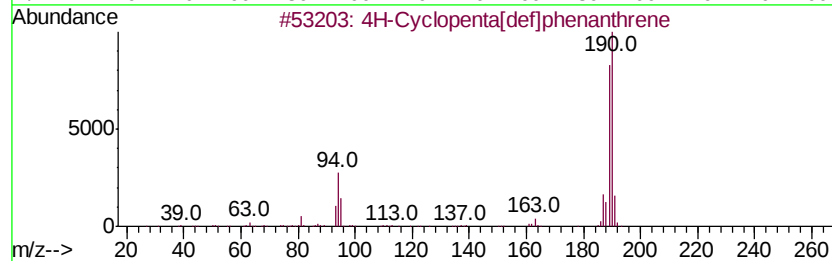
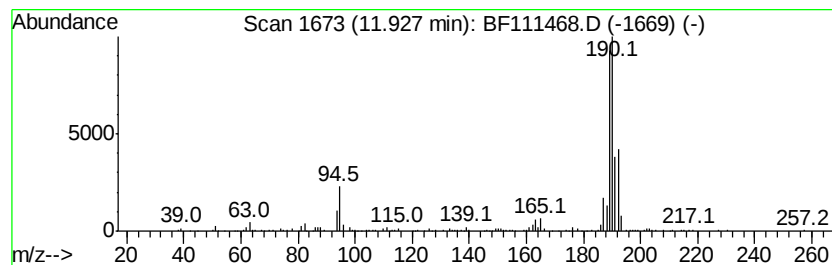
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 4H-Cyclopenta[def]phenanthrene Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.93 | 6.25 ng | 350086 | Phenanthrene-d10 | 11.31 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10     | 000203-64-5  | 76   |
| 2    |    | Thiophene-3-carboxaldehyde, 5-ch... | 190 | C5H4BClO3S | 036155-87-0  | 53   |
| 3    |    | Carbonic acid, ethyl 2-formyl-4,... | 262 | C10H8Cl2O4 | 1000331-34-4 | 50   |
| 4    |    | Methyl diselenide                   | 190 | C2H6Se2    | 007101-31-7  | 43   |
| 5    |    | 2,2'-Bis(4,5-dimethylimidazole)     | 190 | C10H14N4   | 069286-06-2  | 42   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
 Data File : BF111468.D  
 Acq On : 15 Dec 2018 15:03  
 Operator : JU/SJ  
 Sample : J6316-01  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

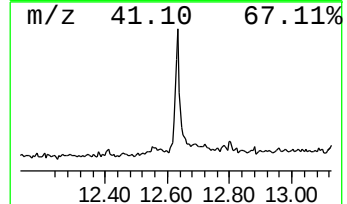
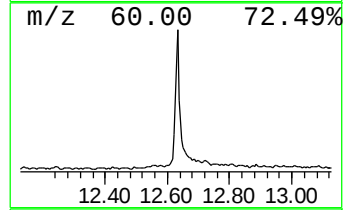
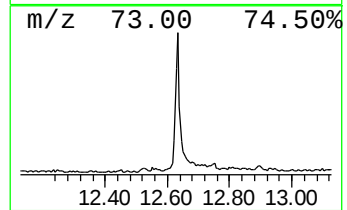
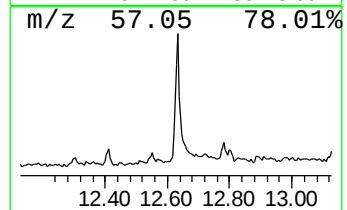
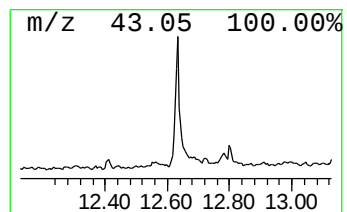
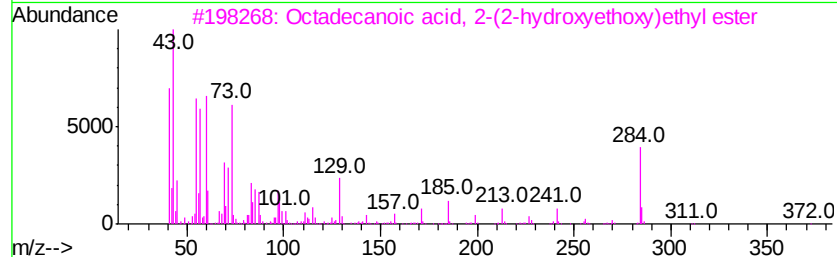
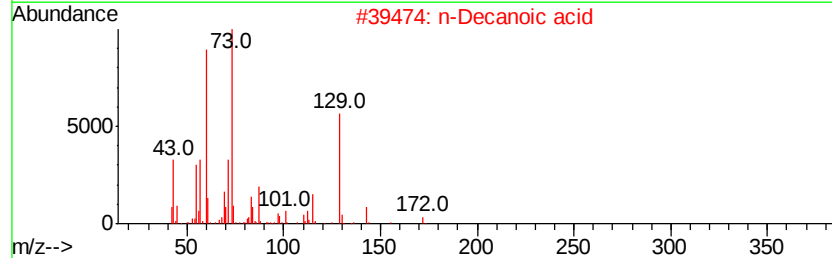
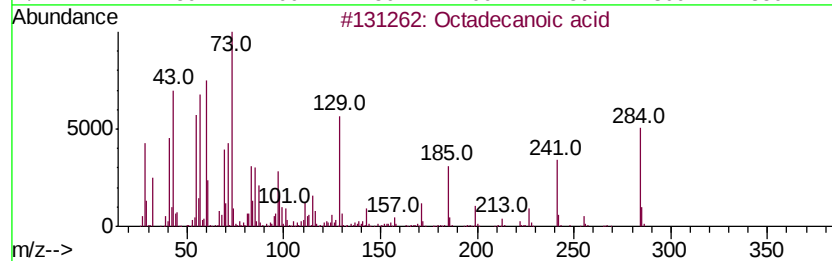
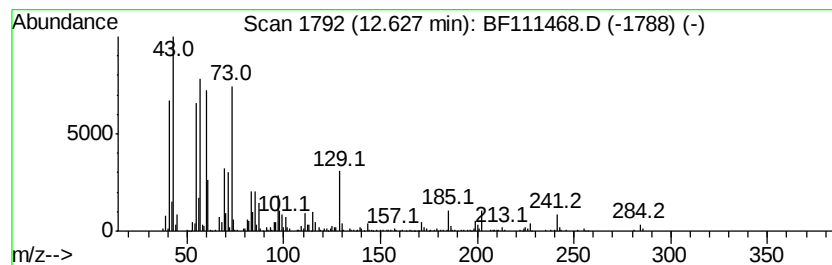
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 17 Octadecanoic acid Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.63 | 8.29 ng | 737116 | Chrysene-d12     | 13.95 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Octadecanoic acid                    | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    | n-Decanoic acid                      | 172 | C10H20O2 | 000334-48-5 | 89   |
| 3    |    | Octadecanoic acid, 2-(2-hydroxyve... | 372 | C22H44O4 | 000106-11-6 | 72   |
| 4    |    | n-Hexadecanoic acid                  | 256 | C16H32O2 | 000057-10-3 | 64   |
| 5    |    | Tridecanoic acid                     | 214 | C13H26O2 | 000638-53-9 | 62   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

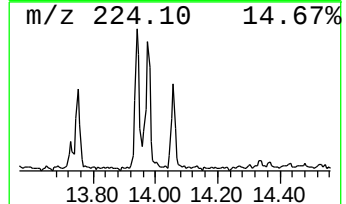
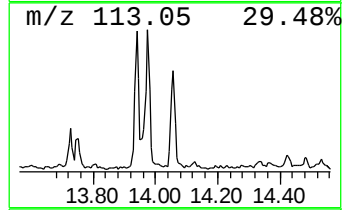
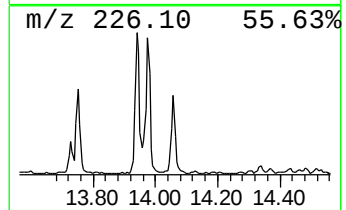
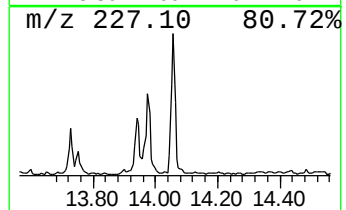
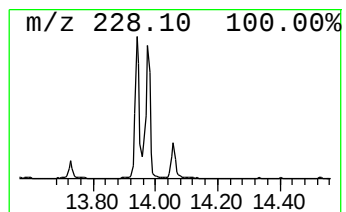
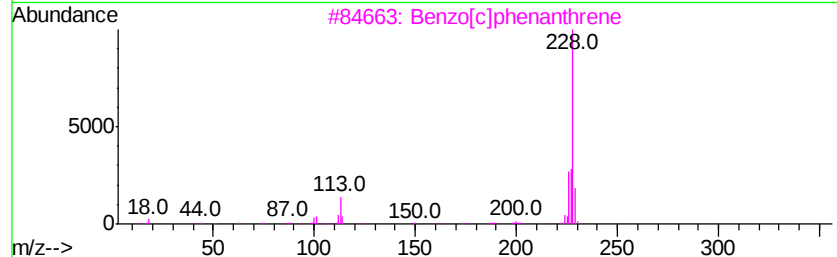
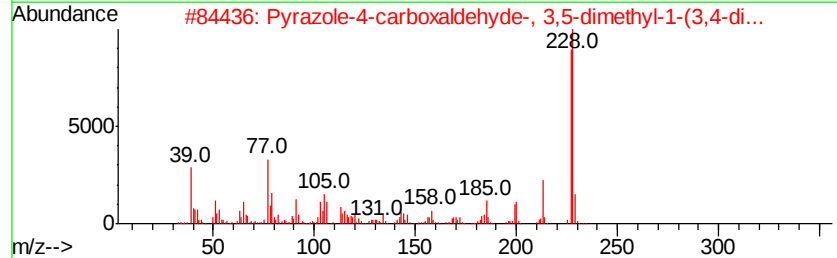
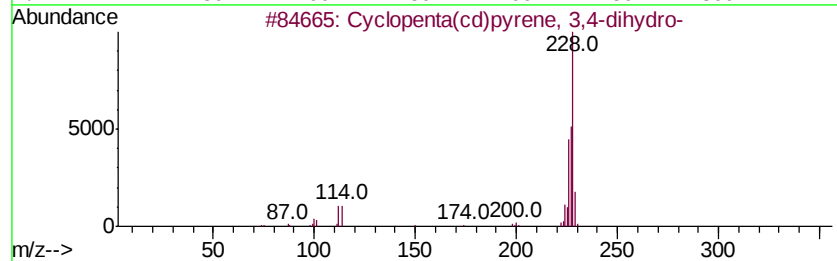
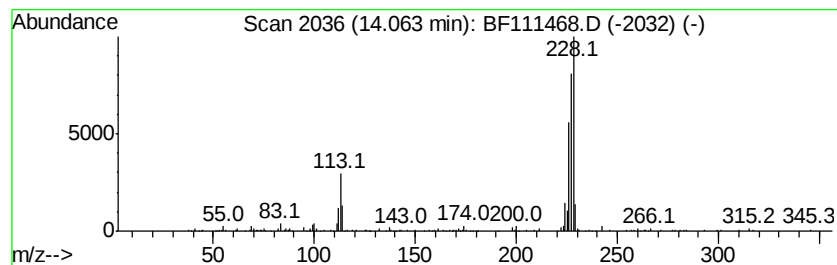
Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.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 Cyclopenta(cd)pyrene, 3,4-d... Concentration Rank 18

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.      |              |      |
|---------|---------|-------------------------------------|------------------|-----------|--------------|------|
| 14.06   | 2.76 ng | 244932                              | Chrysene-d12     | 13.95     |              |      |
| Hit# of | 5       | Tentative ID                        | MW               | MolForm   | CAS#         | Qual |
| 1       |         | Cvclopenta(cd)pvrene, 3,4-dihydro-  | 228              | C18H12    | 025732-74-5  | 78   |
| 2       |         | Pyrazole-4-carboxaldehyde-, 3,5-... | 228              | C14H16N2O | 1000272-54-8 | 50   |
| 3       |         | Benzo[c]lphenanthrene               | 228              | C18H12    | 000195-19-7  | 49   |
| 4       |         | 1,3,5-Triazine-2(1H)-thione, 4-(... | 227              | C9H17N5S  | 023613-02-7  | 46   |
| 5       |         | Dimethyl-[4-[2-(3-methylisoxazol... | 228              | C14H16N2O | 1000306-39-6 | 40   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\  
Data File : BF111468.D  
Acq On : 15 Dec 2018 15:03  
Operator : JU/SJ  
Sample : J6316-01  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

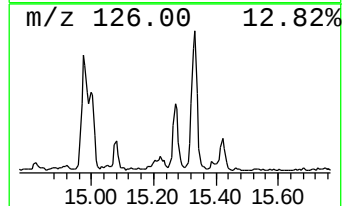
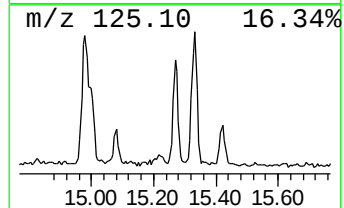
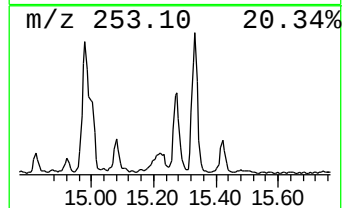
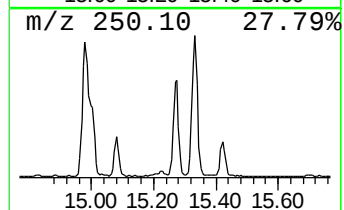
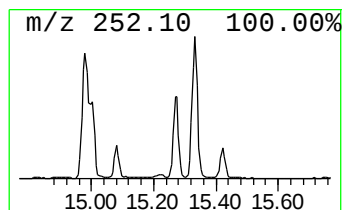
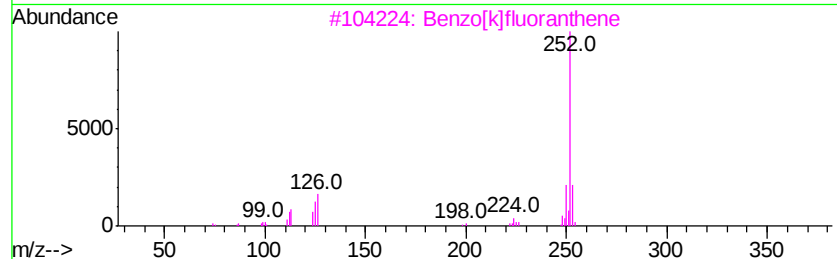
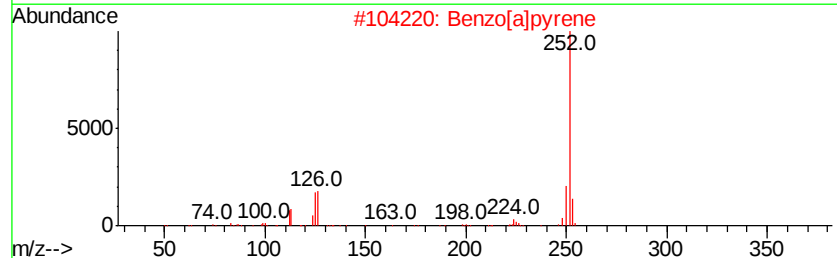
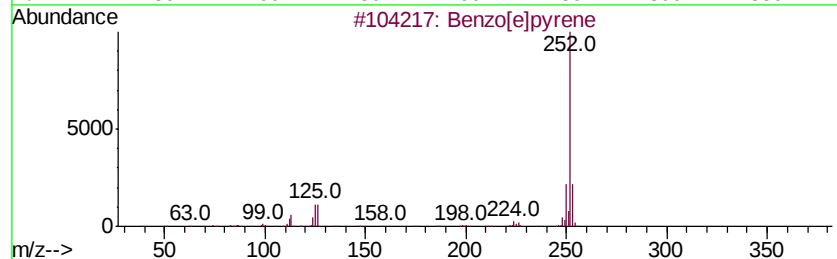
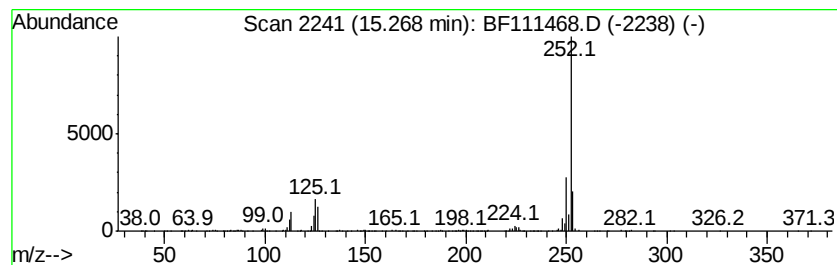
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Benzo[e]pyrene Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.27 | 12.66 ng | 541342 | Perylene-d12     | 15.39 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[e]pyrene            | 252 | C20H12  | 000192-97-2 | 98   |
| 2    |    | Benzo[a]pyrene            | 252 | C20H12  | 000050-32-8 | 96   |
| 3    |    | Benzo[k]fluoranthene      | 252 | C20H12  | 000207-08-9 | 94   |
| 4    |    | Benzo[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 91   |
| 5    |    | Perylene                  | 252 | C20H12  | 000198-55-0 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF121518\  
 Data File : BF111468.D  
 Acq On : 15 Dec 2018 15:03  
 Operator : JU/SJ  
 Sample : J6316-01  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 BORING-SAMPLE-SIGNAL-TOWER-

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF121418.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| 2-Pentanone, 4-hy... | 5.00  | 3.8     | ng    | 199310   | 1                     | 6.79  | 1053370 | 20.0 |
| unknown6.57          | 6.57  | 73.9    | ng    | 3894120  | 1                     | 6.79  | 1053370 | 20.0 |
| Benzene, 1-etheny... | 6.63  | 7.1     | ng    | 375770   | 1                     | 6.79  | 1053370 | 20.0 |
| Benzene, 2-propenyl- | 6.67  | 3.1     | ng    | 164088   | 1                     | 6.79  | 1053370 | 20.0 |
| Indene               | 7.06  | 9.6     | ng    | 507757   | 1                     | 6.79  | 1053370 | 20.0 |
| Benzene, 1-methyl... | 7.41  | 2.7     | ng    | 142215   | 1                     | 6.79  | 1053370 | 20.0 |
| Benzene, (1,3-dim... | 7.65  | 12.2    | ng    | 845890   | 2                     | 8.07  | 1384940 | 20.0 |
| Benzene, 1-methyl... | 8.29  | 8.8     | ng    | 606983   | 2                     | 8.07  | 1384940 | 20.0 |
| Benzene, (1-bromo... | 9.27  | 3.1     | ng    | 226167   | 3                     | 9.83  | 1447950 | 20.0 |
| Anthracene, 2-met... | 11.82 | 3.3     | ng    | 183435   | 4                     | 11.31 | 1120570 | 20.0 |
| n-Hexadecanoic acid  | 11.85 | 18.1    | ng    | 1011690  | 4                     | 11.31 | 1120570 | 20.0 |
| Anthracene, 1-met... | 11.89 | 2.5     | ng    | 140143   | 4                     | 11.31 | 1120570 | 20.0 |
| 4H-Cyclopenta[def... | 11.93 | 6.3     | ng    | 350086   | 4                     | 11.31 | 1120570 | 20.0 |
| Octadecanoic acid    | 12.63 | 8.3     | ng    | 737116   | 5                     | 13.95 | 1777310 | 20.0 |
| Cyclopenta(cd)pyr... | 14.06 | 2.8     | ng    | 244932   | 5                     | 13.95 | 1777310 | 20.0 |
| Benzo[e]pyrene       | 15.27 | 12.7    | ng    | 541342   | 6                     | 15.39 | 855333  | 20.0 |