

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
 Data File : BM002113.D  
 Acq On : 29 Jul 2015 03:28  
 Operator : TP/UM  
 Sample : G3048-13  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C01Y6

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 0.1 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM072715.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 2.763       | 9             | 13          | 17           | rVV      | 12109          | 15121         | 1.00%           | 0.106%        |
| 2         | 2.804       | 17            | 20          | 25           | rVB2     | 5310           | 6851          | 0.45%           | 0.048%        |
| 3         | 3.587       | 149           | 153         | 159          | rVB      | 10533          | 12776         | 0.84%           | 0.089%        |
| 4         | 3.798       | 185           | 189         | 195          | rVB      | 9920           | 13564         | 0.89%           | 0.095%        |
| 5         | 5.239       | 430           | 434         | 436          | rBV      | 8624           | 10592         | 0.70%           | 0.074%        |
| 6         | 5.610       | 491           | 497         | 501          | rBB3     | 7516           | 12209         | 0.81%           | 0.085%        |
| 7         | 6.739       | 683           | 689         | 692          | rBV4     | 7284           | 13908         | 0.92%           | 0.097%        |
| 8         | 6.781       | 692           | 696         | 706          | rVB      | 83080          | 134897        | 8.90%           | 0.943%        |
| 9         | 6.939       | 714           | 723         | 734          | rBV      | 64955          | 108205        | 7.14%           | 0.757%        |
| 10        | 7.069       | 740           | 745         | 750          | rBV3     | 10893          | 17307         | 1.14%           | 0.121%        |
| 11        | 7.133       | 750           | 756         | 766          | rVV      | 93835          | 145904        | 9.62%           | 1.020%        |
| 12        | 7.239       | 771           | 774         | 780          | rVB2     | 7420           | 10274         | 0.68%           | 0.072%        |
| 13        | 7.598       | 823           | 835         | 842          | rVV      | 232956         | 370396        | 24.43%          | 2.590%        |
| 14        | 7.710       | 849           | 854         | 858          | rVV3     | 3966           | 7100          | 0.47%           | 0.050%        |
| 15        | 8.133       | 923           | 926         | 931          | rVV4     | 6681           | 9148          | 0.60%           | 0.064%        |
| 16        | 8.233       | 936           | 943         | 949          | rVV4     | 5062           | 11199         | 0.74%           | 0.078%        |
| 17        | 8.316       | 952           | 957         | 966          | rBV      | 62463          | 103657        | 6.84%           | 0.725%        |
| 18        | 8.416       | 966           | 974         | 980          | rVV6     | 14052          | 29695         | 1.96%           | 0.208%        |
| 19        | 8.592       | 998           | 1004        | 1008         | rVV4     | 5409           | 8746          | 0.58%           | 0.061%        |
| 20        | 8.686       | 1015          | 1020        | 1025         | rVV6     | 7881           | 14970         | 0.99%           | 0.105%        |
| 21        | 8.757       | 1025          | 1032        | 1038         | rVV      | 82010          | 134777        | 8.89%           | 0.942%        |
| 22        | 8.816       | 1038          | 1042        | 1047         | rVV2     | 8694           | 14351         | 0.95%           | 0.100%        |
| 23        | 8.910       | 1050          | 1058        | 1059         | rBV5     | 4337           | 8933          | 0.59%           | 0.062%        |
| 24        | 8.933       | 1059          | 1062        | 1068         | rVB5     | 5455           | 10563         | 0.70%           | 0.074%        |
| 25        | 9.222       | 1107          | 1111        | 1116         | rBV4     | 4709           | 9257          | 0.61%           | 0.065%        |
| 26        | 9.298       | 1119          | 1124        | 1130         | rVB4     | 7185           | 12760         | 0.84%           | 0.089%        |
| 27        | 9.475       | 1145          | 1154        | 1163         | rBV2     | 51909          | 93733         | 6.18%           | 0.655%        |
| 28        | 9.557       | 1163          | 1168        | 1172         | rBV3     | 9674           | 17644         | 1.16%           | 0.123%        |
| 29        | 9.727       | 1191          | 1197        | 1202         | rVB5     | 3934           | 8096          | 0.53%           | 0.057%        |
| 30        | 9.792       | 1202          | 1208        | 1213         | rBV5     | 5916           | 13971         | 0.92%           | 0.098%        |
| 31        | 10.016      | 1241          | 1246        | 1254         | rVB      | 105160         | 166847        | 11.00%          | 1.167%        |
| 32        | 10.169      | 1266          | 1272        | 1276         | rBV3     | 6887           | 11758         | 0.78%           | 0.082%        |
| 33        | 10.251      | 1281          | 1286        | 1290         | rBV2     | 12933          | 19591         | 1.29%           | 0.137%        |
| 34        | 10.386      | 1298          | 1309        | 1315         | rBV      | 309754         | 534364        | 35.24%          | 3.736%        |

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Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM072715.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 10.433 | 1315 | 1317 | 1322 | rVB  | 15104  | 20419  | 1.35%  | 0.143% |
| 36 | 10.533 | 1330 | 1334 | 1343 | rVB2 | 48844  | 83313  | 5.49%  | 0.582% |
| 37 | 10.774 | 1371 | 1375 | 1381 | rVB6 | 3967   | 6690   | 0.44%  | 0.047% |
| 38 | 10.863 | 1384 | 1390 | 1397 | rVB2 | 7516   | 13856  | 0.91%  | 0.097% |
| 39 | 10.927 | 1397 | 1401 | 1404 | rBV4 | 6808   | 11679  | 0.77%  | 0.082% |
| 40 | 10.992 | 1407 | 1412 | 1415 | rBV3 | 5122   | 7529   | 0.50%  | 0.053% |
| 41 | 11.069 | 1423 | 1425 | 1430 | rVB3 | 6006   | 9626   | 0.63%  | 0.067% |
| 42 | 11.139 | 1435 | 1437 | 1442 | rVV2 | 4920   | 7878   | 0.52%  | 0.055% |
| 43 | 11.210 | 1444 | 1449 | 1457 | rVV6 | 6530   | 16734  | 1.10%  | 0.117% |
| 44 | 11.292 | 1459 | 1463 | 1471 | rVB4 | 8826   | 16737  | 1.10%  | 0.117% |
| 45 | 11.404 | 1477 | 1482 | 1490 | rVV4 | 24754  | 53835  | 3.55%  | 0.376% |
| 46 | 11.469 | 1491 | 1493 | 1498 | rVV5 | 8860   | 14393  | 0.95%  | 0.101% |
| 47 | 11.569 | 1502 | 1510 | 1514 | rBV5 | 5153   | 10754  | 0.71%  | 0.075% |
| 48 | 11.663 | 1522 | 1526 | 1531 | rVB5 | 8970   | 14016  | 0.92%  | 0.098% |
| 49 | 11.810 | 1544 | 1551 | 1553 | rBV6 | 11127  | 18513  | 1.22%  | 0.129% |
| 50 | 12.057 | 1589 | 1593 | 1600 | rVB  | 22183  | 38182  | 2.52%  | 0.267% |
| 51 | 12.121 | 1600 | 1604 | 1611 | rBV6 | 8160   | 15163  | 1.00%  | 0.106% |
| 52 | 12.210 | 1616 | 1619 | 1624 | rVB6 | 5602   | 8702   | 0.57%  | 0.061% |
| 53 | 12.280 | 1624 | 1631 | 1635 | rBV3 | 10654  | 23120  | 1.52%  | 0.162% |
| 54 | 12.451 | 1655 | 1660 | 1664 | rBV4 | 9941   | 19214  | 1.27%  | 0.134% |
| 55 | 12.592 | 1681 | 1684 | 1689 | rBV6 | 10503  | 16191  | 1.07%  | 0.113% |
| 56 | 12.674 | 1695 | 1698 | 1703 | rVB6 | 10850  | 16270  | 1.07%  | 0.114% |
| 57 | 12.727 | 1704 | 1707 | 1712 | rVV7 | 8993   | 16384  | 1.08%  | 0.115% |
| 58 | 12.780 | 1712 | 1716 | 1722 | rVB  | 40382  | 56976  | 3.76%  | 0.398% |
| 59 | 12.851 | 1722 | 1728 | 1735 | rBV8 | 18992  | 41719  | 2.75%  | 0.292% |
| 60 | 13.021 | 1753 | 1757 | 1760 | rBV2 | 25344  | 36643  | 2.42%  | 0.256% |
| 61 | 13.074 | 1764 | 1766 | 1772 | rVB5 | 12797  | 23666  | 1.56%  | 0.165% |
| 62 | 13.415 | 1821 | 1824 | 1829 | rBV3 | 12250  | 24860  | 1.64%  | 0.174% |
| 63 | 13.580 | 1848 | 1852 | 1857 | rVB7 | 14589  | 23218  | 1.53%  | 0.162% |
| 64 | 13.674 | 1862 | 1868 | 1872 | rVV  | 242305 | 345276 | 22.77% | 2.414% |
| 65 | 13.721 | 1872 | 1876 | 1883 | rVB2 | 114930 | 217709 | 14.36% | 1.522% |
| 66 | 13.780 | 1883 | 1886 | 1891 | rBV6 | 17360  | 26832  | 1.77%  | 0.188% |
| 67 | 13.951 | 1910 | 1915 | 1920 | rBV  | 202947 | 294034 | 19.39% | 2.056% |
| 68 | 14.080 | 1933 | 1937 | 1944 | rVB2 | 65825  | 101556 | 6.70%  | 0.710% |
| 69 | 14.157 | 1945 | 1950 | 1955 | rBV4 | 41241  | 64928  | 4.28%  | 0.454% |
| 70 | 14.221 | 1956 | 1961 | 1962 | rVV2 | 30361  | 46547  | 3.07%  | 0.325% |
| 71 | 14.262 | 1963 | 1968 | 1974 | rVB2 | 462347 | 715103 | 47.16% | 5.000% |

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Peak Location: TOP

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Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM072715.M  
Title : SVOA CALIBRATION

|     |        |      |      |      |      |        |         |         |         |
|-----|--------|------|------|------|------|--------|---------|---------|---------|
| 72  | 14.492 | 2003 | 2007 | 2012 | rVB2 | 48896  | 77376   | 5.10%   | 0.541%  |
| 73  | 14.662 | 2033 | 2036 | 2040 | rVB  | 28504  | 35176   | 2.32%   | 0.246%  |
| 74  | 14.780 | 2053 | 2056 | 2063 | rVB3 | 36820  | 52641   | 3.47%   | 0.368%  |
| 75  | 14.886 | 2070 | 2074 | 2076 | rBV4 | 25463  | 38656   | 2.55%   | 0.270%  |
| 76  | 14.968 | 2085 | 2088 | 2096 | rBV7 | 24014  | 46784   | 3.09%   | 0.327%  |
| 77  | 15.045 | 2097 | 2101 | 2105 | rVV4 | 43530  | 65020   | 4.29%   | 0.455%  |
| 78  | 15.257 | 2132 | 2137 | 2142 | rVB  | 233306 | 325642  | 21.48%  | 2.277%  |
| 79  | 15.521 | 2177 | 2182 | 2186 | rVB  | 97507  | 142234  | 9.38%   | 0.994%  |
| 80  | 15.998 | 2256 | 2263 | 2268 | rBV  | 261811 | 454713  | 29.99%  | 3.179%  |
| 81  | 16.821 | 2399 | 2403 | 2407 | rVB2 | 165207 | 217002  | 14.31%  | 1.517%  |
| 82  | 17.009 | 2430 | 2435 | 2440 | rBV  | 592131 | 880803  | 58.09%  | 6.158%  |
| 83  | 17.051 | 2440 | 2442 | 2447 | rVV  | 111004 | 141668  | 9.34%   | 0.990%  |
| 84  | 17.109 | 2448 | 2452 | 2456 | rVV  | 251956 | 333920  | 22.02%  | 2.335%  |
| 85  | 17.421 | 2502 | 2505 | 2511 | rBV4 | 66550  | 96647   | 6.37%   | 0.676%  |
| 86  | 17.909 | 2586 | 2588 | 2591 | rBV2 | 69124  | 56913   | 3.75%   | 0.398%  |
| 87  | 18.080 | 2614 | 2617 | 2622 | rVB2 | 134082 | 172906  | 11.40%  | 1.209%  |
| 88  | 18.633 | 2708 | 2711 | 2715 | rVB2 | 138628 | 155076  | 10.23%  | 1.084%  |
| 89  | 19.080 | 2783 | 2787 | 2792 | rBV  | 241736 | 323921  | 21.36%  | 2.265%  |
| 90  | 19.133 | 2792 | 2796 | 2802 | rVV  | 335319 | 444865  | 29.34%  | 3.110%  |
| 91  | 19.221 | 2808 | 2811 | 2815 | rVB  | 289351 | 345673  | 22.80%  | 2.417%  |
| 92  | 19.415 | 2840 | 2844 | 2847 | rBV  | 350376 | 525215  | 34.64%  | 3.672%  |
| 93  | 19.668 | 2885 | 2887 | 2894 | rVB2 | 241121 | 314117  | 20.72%  | 2.196%  |
| 94  | 19.933 | 2929 | 2932 | 2938 | rBV5 | 230015 | 289886  | 19.12%  | 2.027%  |
| 95  | 21.127 | 3132 | 3135 | 3139 | rVB  | 262345 | 297916  | 19.65%  | 2.083%  |
| 96  | 21.209 | 3145 | 3149 | 3161 | rVB2 | 885676 | 1516272 | 100.00% | 10.601% |
| 97  | 21.785 | 3244 | 3247 | 3253 | rVB  | 320922 | 363298  | 23.96%  | 2.540%  |
| 98  | 22.238 | 3321 | 3324 | 3328 | rVB  | 409579 | 521020  | 34.36%  | 3.643%  |
| 99  | 23.279 | 3498 | 3501 | 3506 | rBV2 | 400877 | 550527  | 36.31%  | 3.849%  |
| 100 | 23.421 | 3521 | 3525 | 3532 | rVB2 | 538307 | 925523  | 61.04%  | 6.471%  |

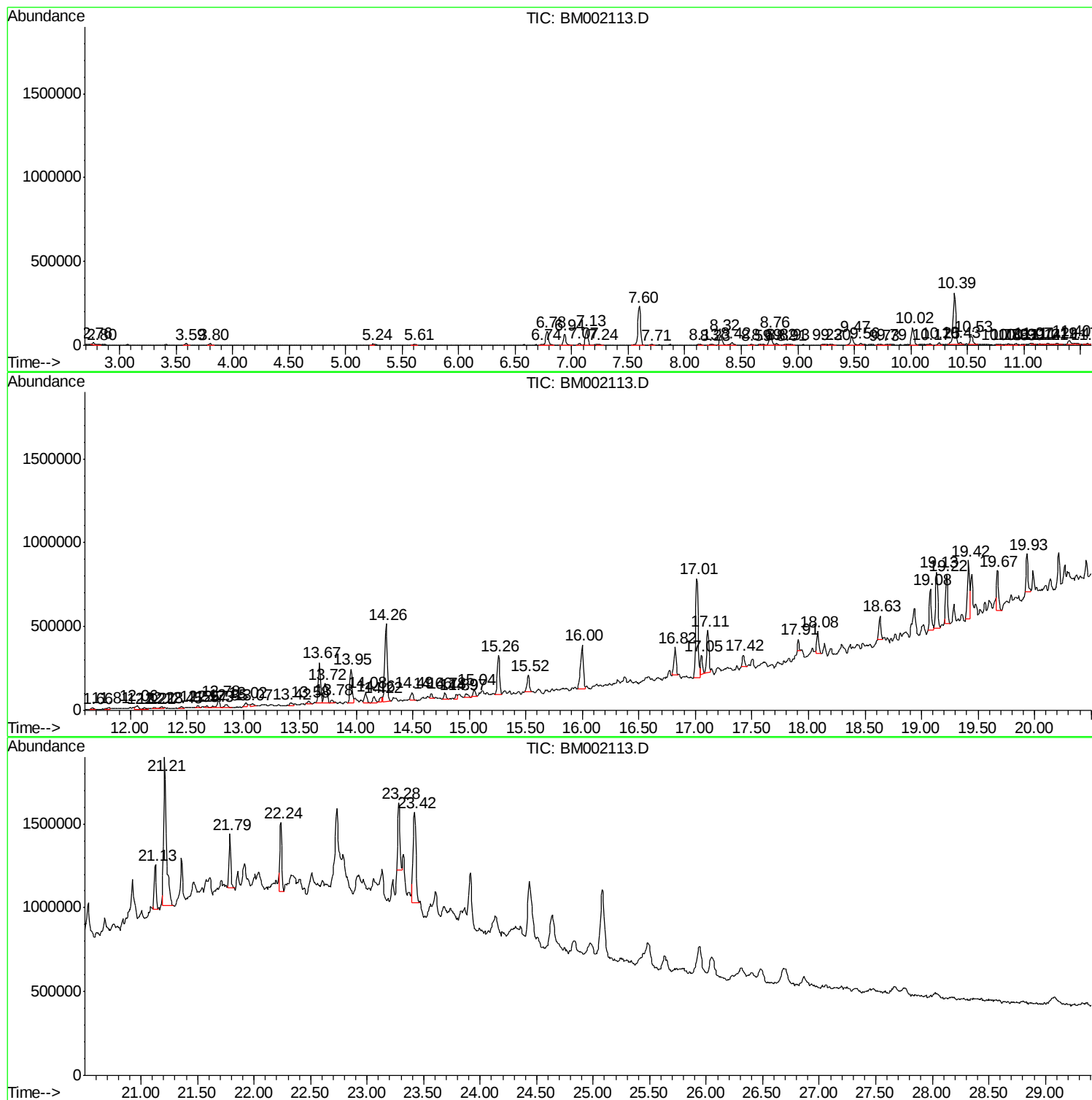
Sum of corrected areas: 14302839

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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

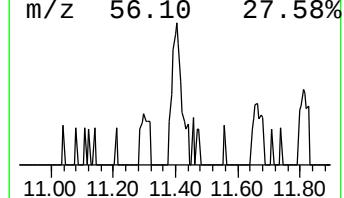
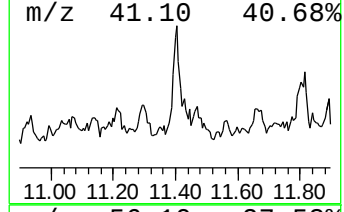
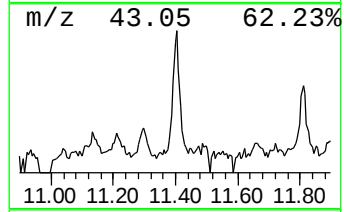
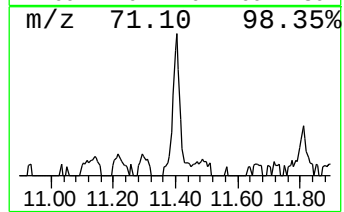
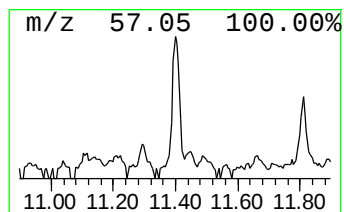
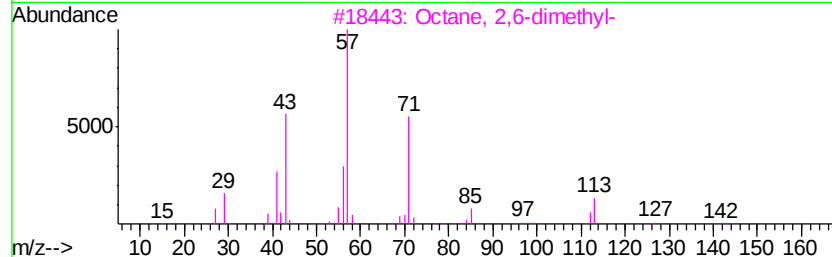
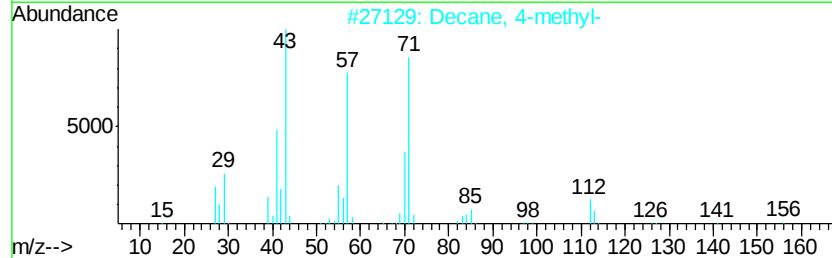
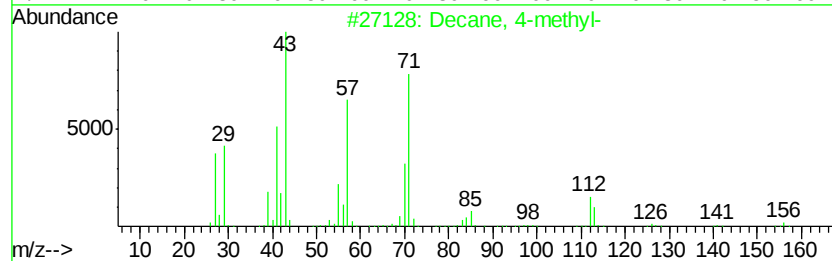
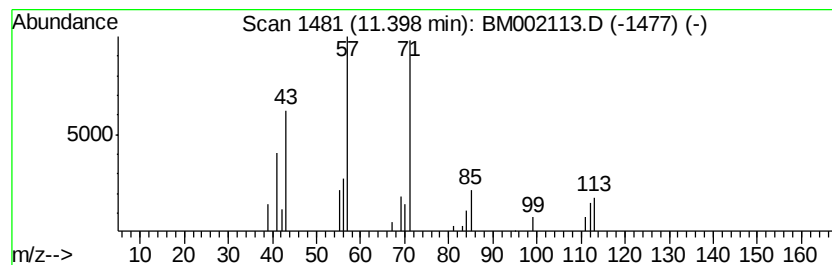
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 11.40 | 2.01 ng/ul | 53835 | Napthalene-d8    | 10.39 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 4-methyl-          | 156 | C11H24  | 002847-72-5 | 64   |
| 2    |    | Decane, 4-methyl-          | 156 | C11H24  | 002847-72-5 | 64   |
| 3    |    | Octane, 2,6-dimethyl-      | 142 | C10H22  | 002051-30-1 | 64   |
| 4    |    | Octane, 3,5-dimethyl-      | 142 | C10H22  | 015869-93-9 | 59   |
| 5    |    | Nonane, 4-methyl-5-propyl- | 184 | C13H28  | 062185-55-1 | 59   |



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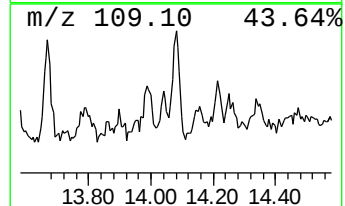
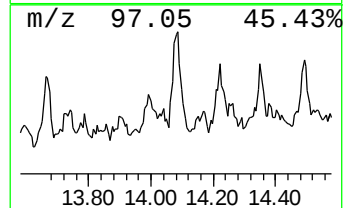
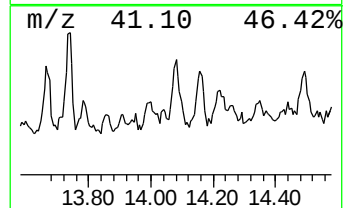
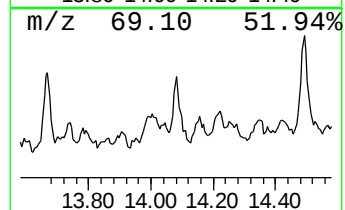
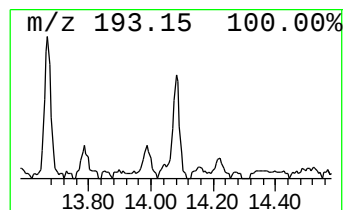
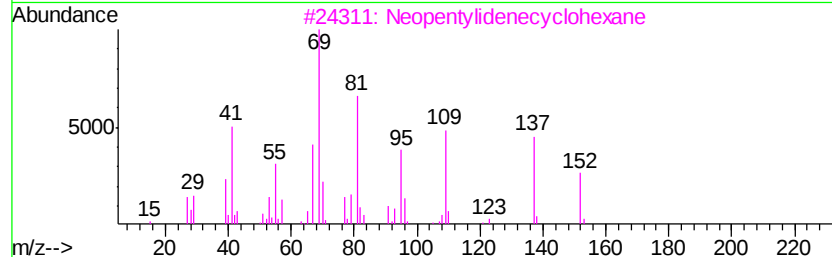
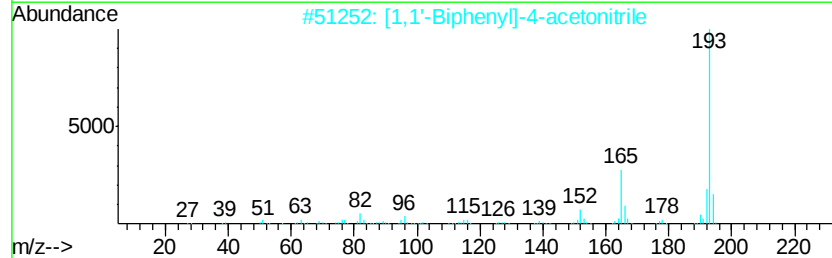
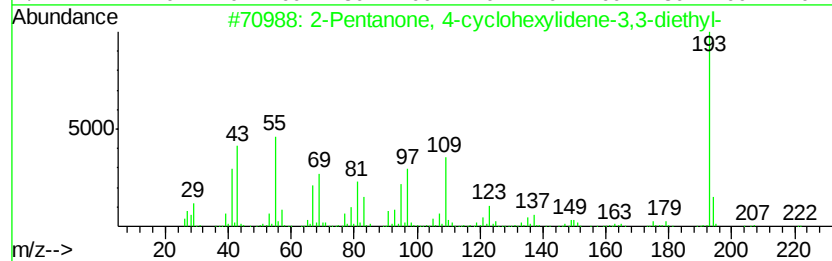
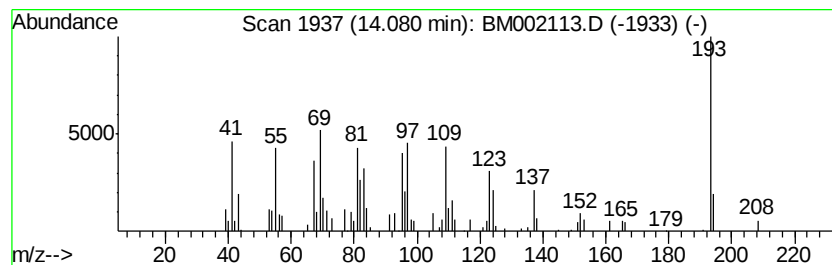
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\*\*\*\*\*  
Peak Number 2 unknown-01 Concentration Rank 12

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.         |      |      |
|---------|-------------------------------------|--------------|------------------|--------------|------|------|
| 14.08   | 2.84 ng/ul                          | 101556       | Acenaphthene-d10 | 14.26        |      |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm      | CAS# | Qual |
| 1       | 2-Pentanone, 4-cvclohexylidene-3... | 222          | C15H26O          | 313253-65-5  | 28   |      |
| 2       | [1,1'-Biphenyl]-4-acetonitrile      | 193          | C14H11N          | 031603-77-7  | 27   |      |
| 3       | Neopentylidenecyclohexane           | 152          | C11H20           | 039546-80-0  | 25   |      |
| 4       | 4H-1,3,2-Dioxaborin, 4-ethenyl-2... | 208          | C12H21BO2        | 074630-04-9  | 23   |      |
| 5       | S-Phenyl 5-(phenylthio)pentaneth... | 302          | C17H18OS2        | 1000234-40-7 | 17   |      |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

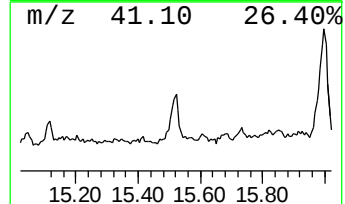
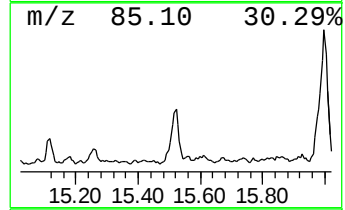
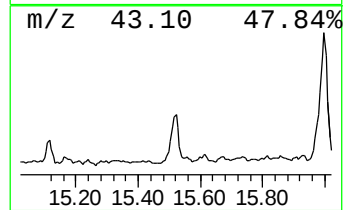
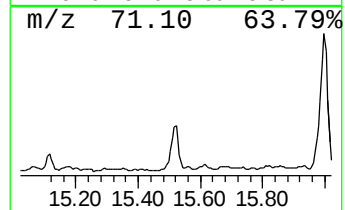
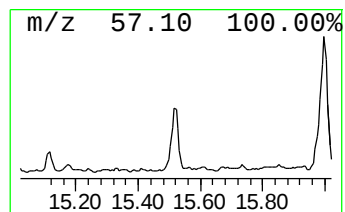
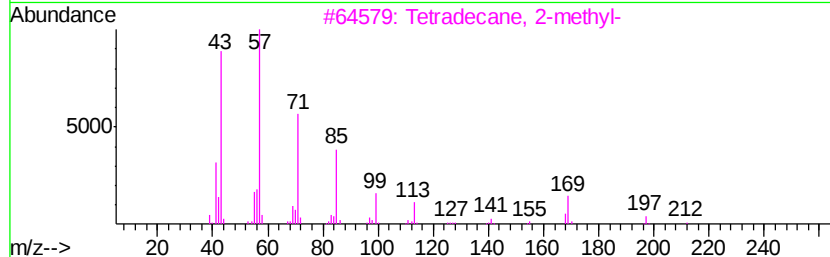
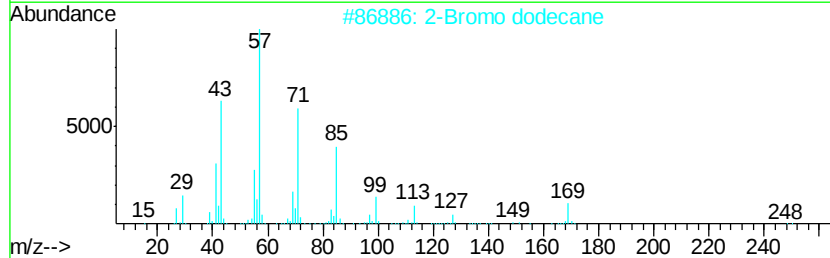
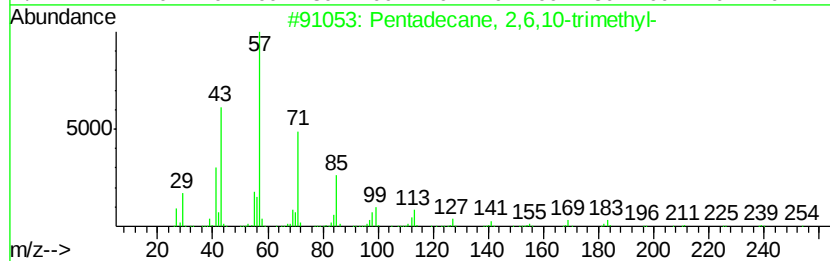
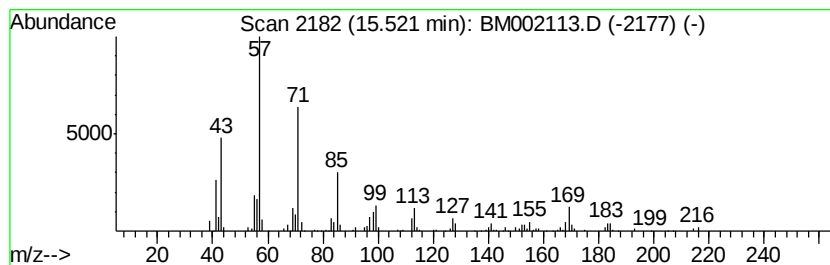
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.52 | 3.98 ng/ul | 142234 | Acenaphthene-d10 | 14.26 |

| Hit# | of | Tentative ID                   | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------|-----|----------|-------------|------|
| 1    | 5  | Pentadecane, 2,6,10-trimethyl- | 254 | C18H38   | 003892-00-0 | 80   |
| 2    |    | 2-Bromo dodecane               | 248 | C12H25Br | 013187-99-0 | 80   |
| 3    |    | Tetradecane, 2-methyl-         | 212 | C15H32   | 001560-95-8 | 80   |
| 4    |    | Tridecane                      | 184 | C13H28   | 000629-50-5 | 74   |
| 5    |    | Eicosane                       | 282 | C20H42   | 000112-95-8 | 64   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

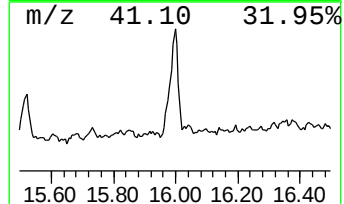
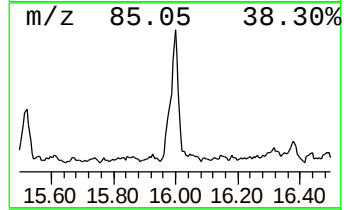
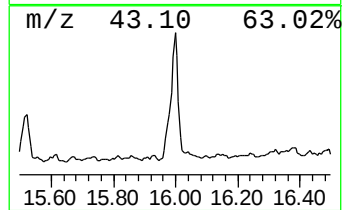
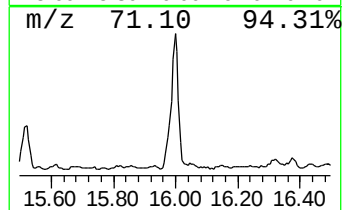
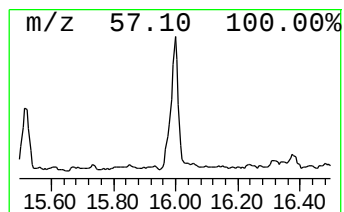
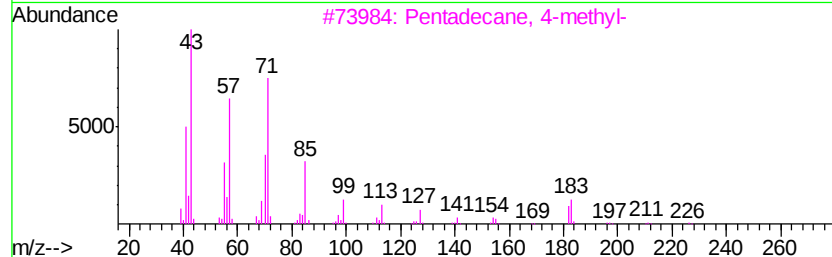
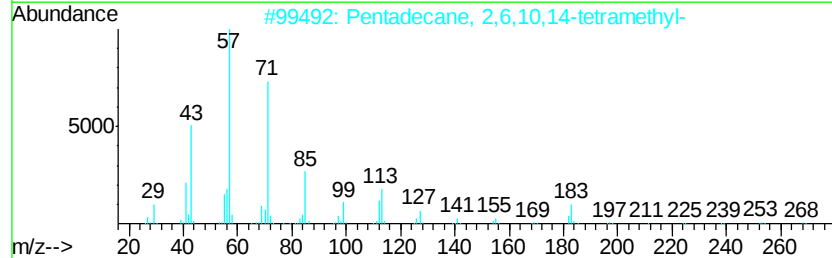
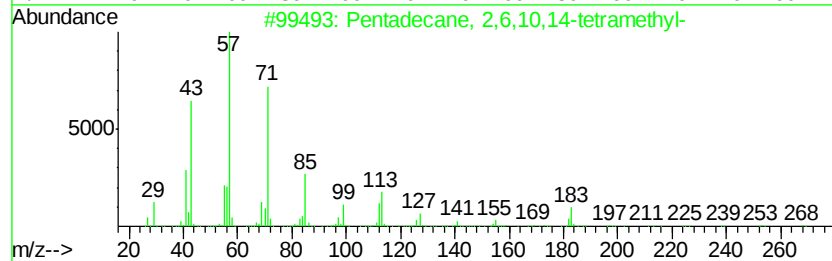
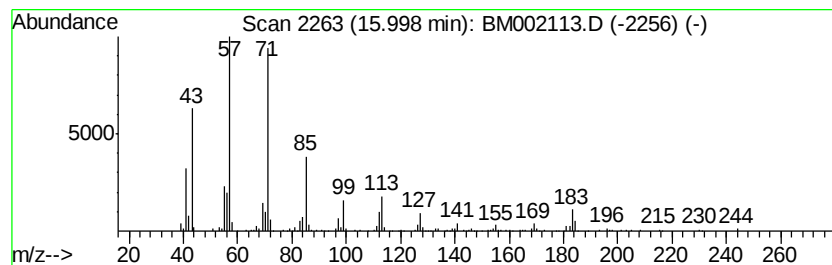
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 16.00 | 10.33 ng/ul | 454713 | Phenanthrene-d10 | 17.01 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 91   |
| 2    |    | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 90   |
| 3    |    | Pentadecane, 4-methyl-              | 226 | C16H34  | 002801-87-8 | 86   |
| 4    |    | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 80   |
| 5    |    | Dodecane, 4,6-dimethyl-             | 198 | C14H30  | 061141-72-8 | 74   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
 Data File : BM002113.D  
 Acq On : 29 Jul 2015 03:28  
 Operator : TP/UM  
 Sample : G3048-13  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
 Quant Title : SVOA CALIBRATION

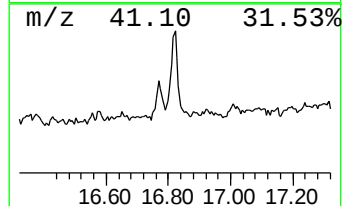
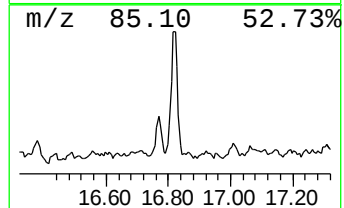
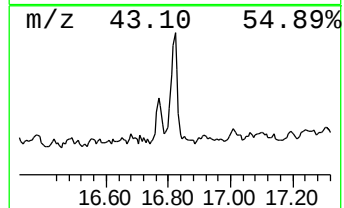
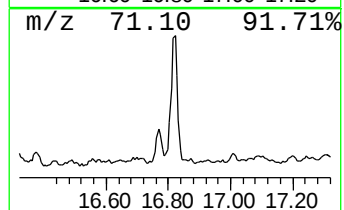
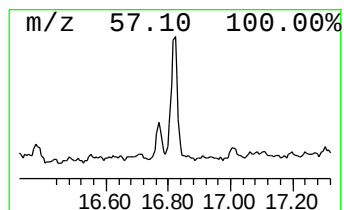
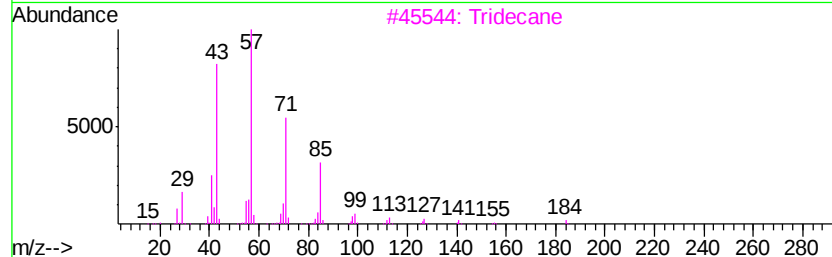
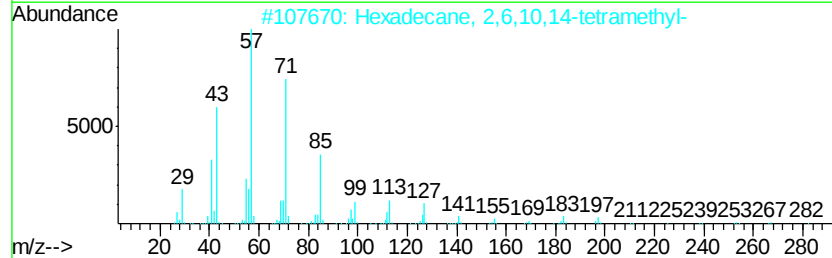
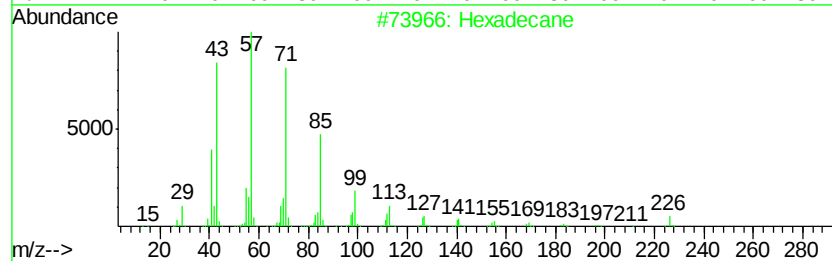
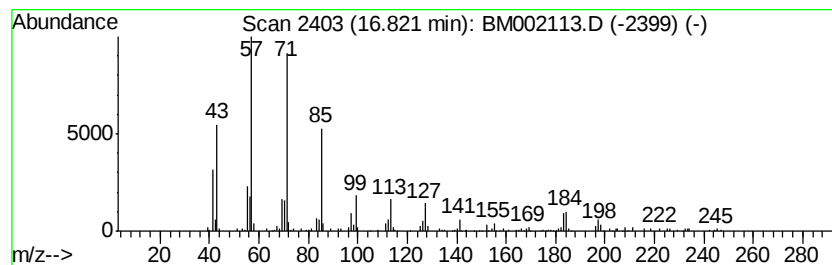
TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.82 | 4.93 ng/ul | 217002 | Phenanthrene-d10 | 17.01 |

| Hit# of | 5                                  | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|------------------------------------|--------------|-----|---------|-------------|------|
| 1       | Hexadecane                         |              | 226 | C16H34  | 000544-76-3 | 91   |
| 2       | Hexadecane, 2,6,10,14-tetramethyl- |              | 282 | C20H42  | 000638-36-8 | 90   |
| 3       | Tridecane                          |              | 184 | C13H28  | 000629-50-5 | 90   |
| 4       | Tetradecane                        |              | 198 | C14H30  | 000629-59-4 | 87   |
| 5       | Pentadecane, 4-methyl-             |              | 226 | C16H34  | 002801-87-8 | 87   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

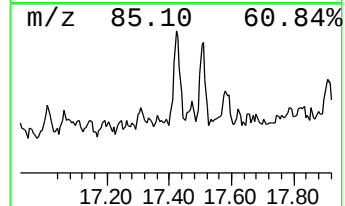
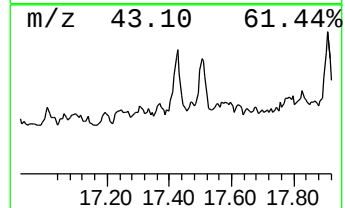
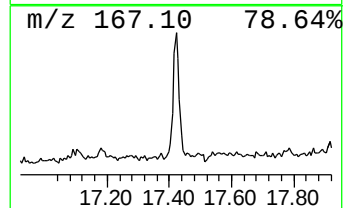
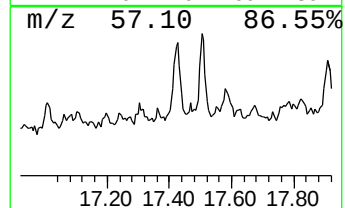
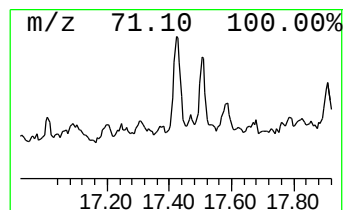
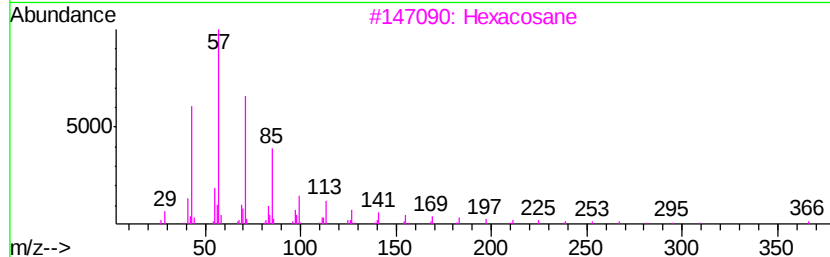
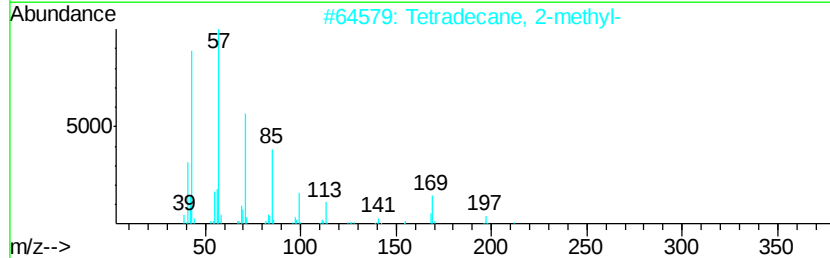
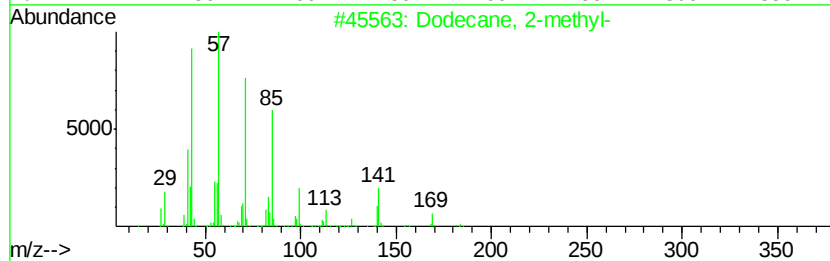
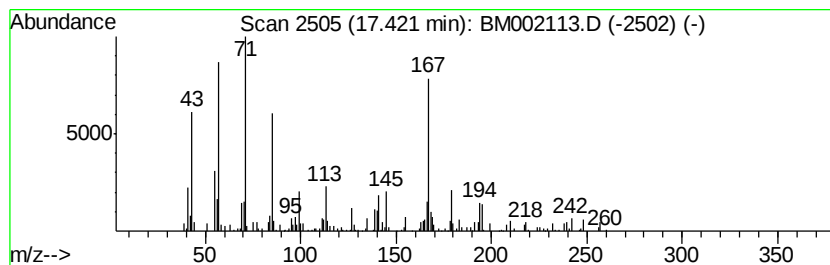
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 17.42 | 2.19 ng/ul | 96647 | Phenanthrene-d10 | 17.01 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane, 2-methyl-     | 184 | C13H28  | 001560-97-0 | 52   |
| 2    |    | Tetradecane, 2-methyl-  | 212 | C15H32  | 001560-95-8 | 45   |
| 3    |    | Hexacosane              | 366 | C26H54  | 000630-01-3 | 43   |
| 4    |    | Undecane, 4,8-dimethyl- | 184 | C13H28  | 017301-33-6 | 38   |
| 5    |    | Tridecane, 3-methyl-    | 198 | C14H30  | 006418-41-3 | 38   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

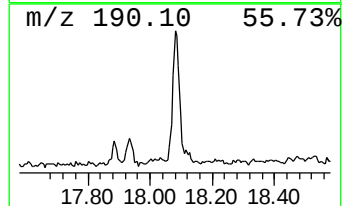
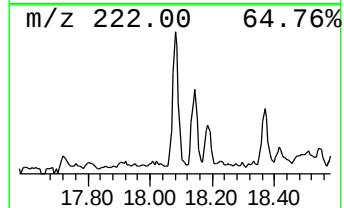
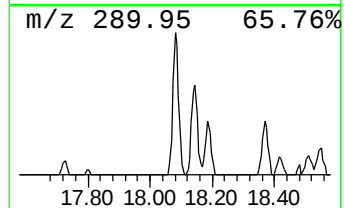
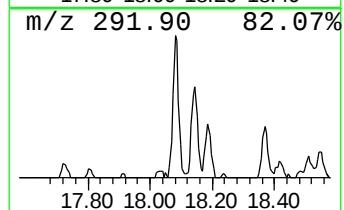
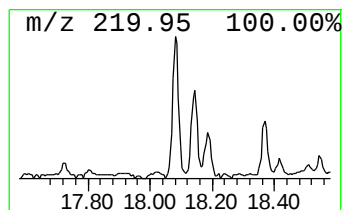
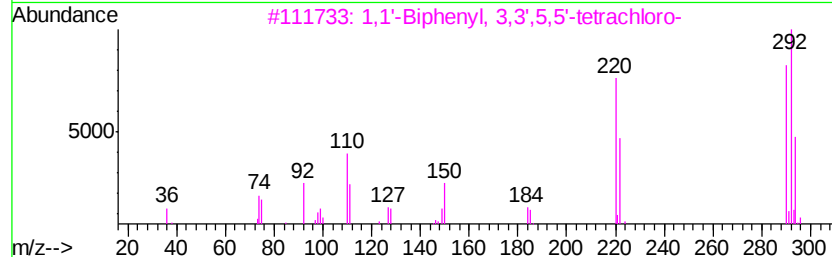
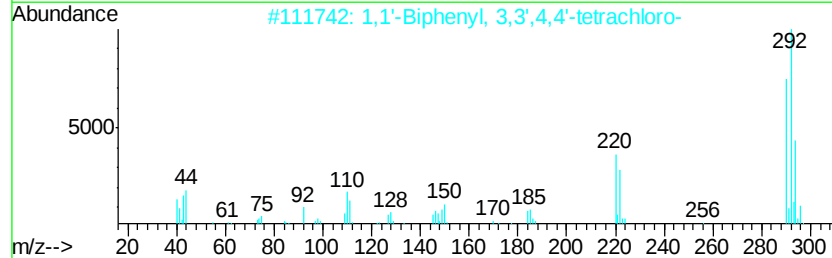
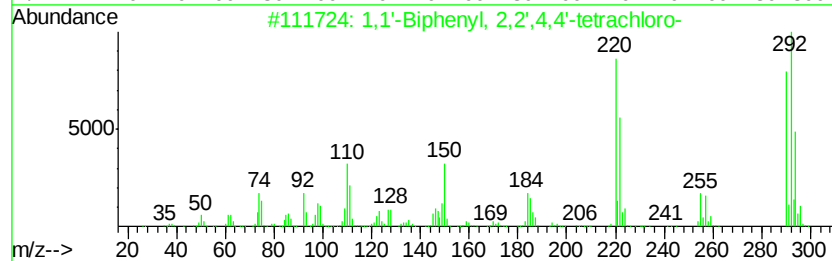
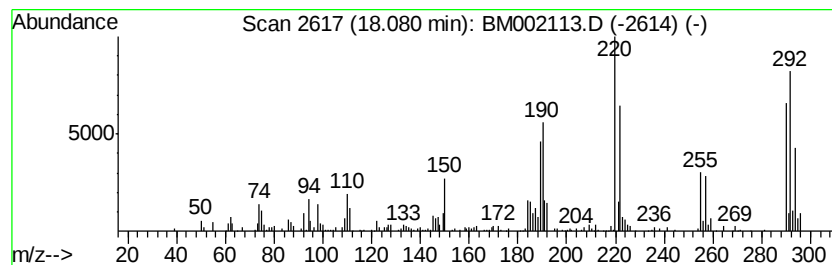
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.08 | 3.93 ng/ul | 172906 | Phenanthrene-d10 | 17.01 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4 | 002437-79-8 | 96   |
| 2    |    |   | 1,1'-Biphenyl, 3,3',4,4'-tetrach... | 290 | C12H6Cl4 | 032598-13-3 | 95   |
| 3    |    |   | 1,1'-Biphenyl, 3,3',5,5'-tetrach... | 290 | C12H6Cl4 | 033284-52-5 | 95   |
| 4    |    |   | 1,1'-Biphenyl, 2,2',5,5'-tetrach... | 290 | C12H6Cl4 | 035693-99-3 | 94   |
| 5    |    |   | 1,1'-Biphenyl, 2,3,4,5-tetrachloro- | 290 | C12H6Cl4 | 033284-53-6 | 93   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

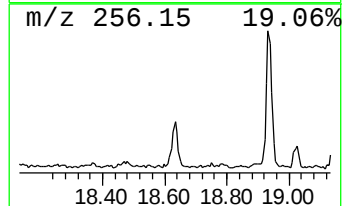
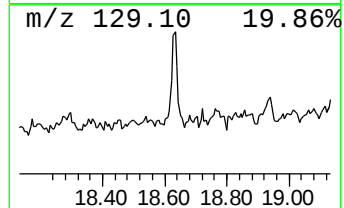
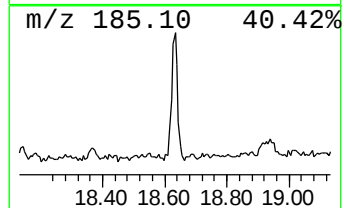
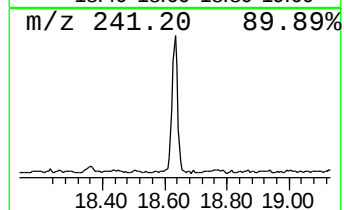
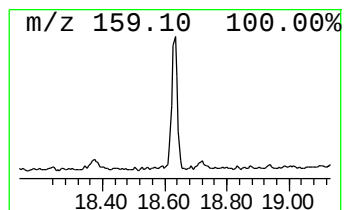
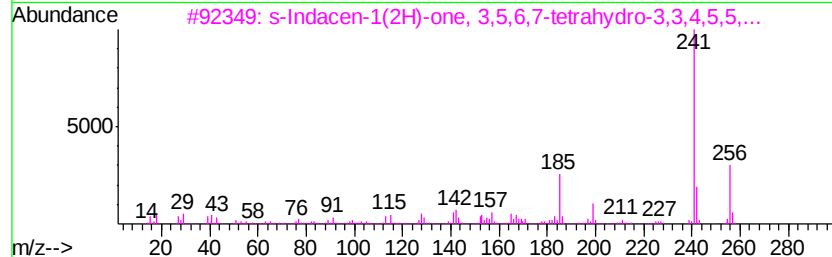
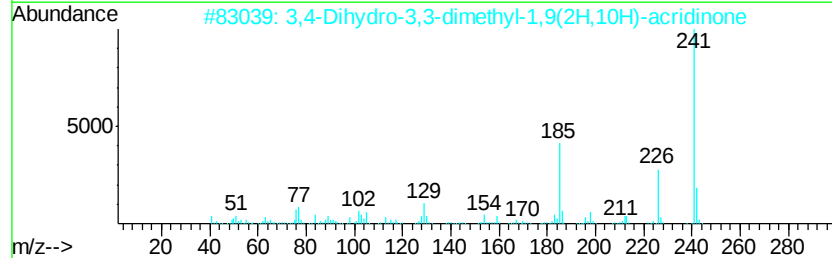
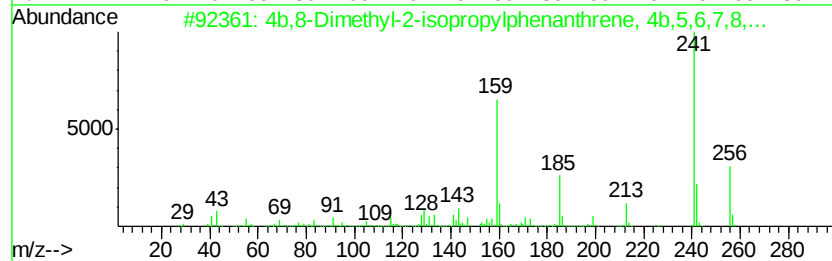
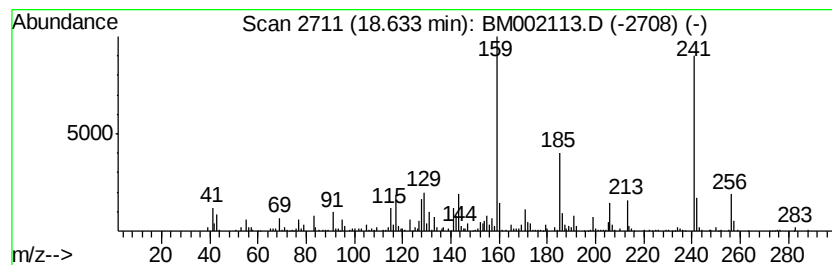
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 4b,8-Dimethyl-2-isopropylph... Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.63 | 3.52 ng/ul | 155076 | Phenanthrene-d10 | 17.01 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 4b,8-Dimethyl-2-isopropylphenant... | 256 | C19H28    | 1000197-14-1 | 95   |
| 2    |    | 3,4-Dihydro-3,3-dimethyl-1,9(2H,... | 241 | C15H15N02 | 1000212-95-2 | 43   |
| 3    |    | s-Indacen-1(2H)-one, 3,5,6,7-tet... | 256 | C18H24O   | 038754-94-8  | 35   |
| 4    |    | As-Indacene, 1,2,3,6,7,8-hexahyd... | 256 | C19H28    | 017465-47-3  | 30   |
| 5    |    | 1-Phenanthrenecarboxaldehyde, 1,... | 284 | C20H28O   | 024035-50-5  | 25   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

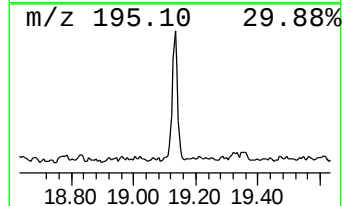
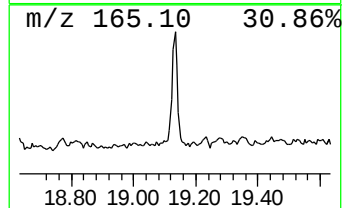
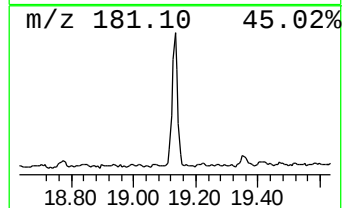
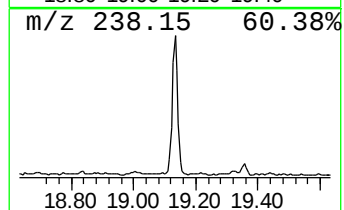
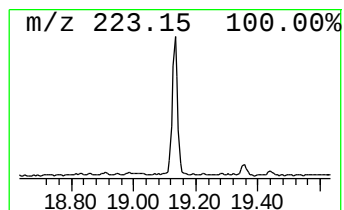
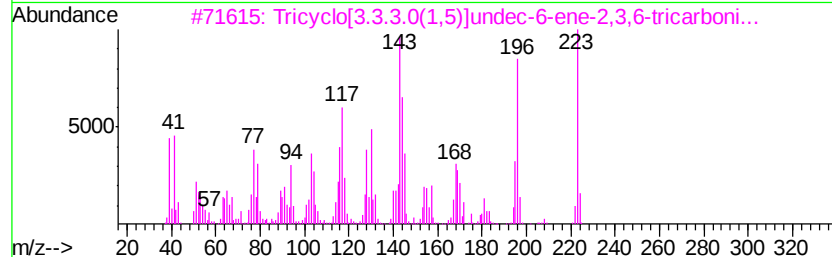
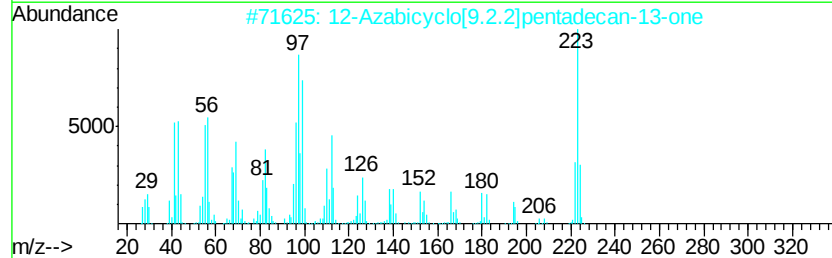
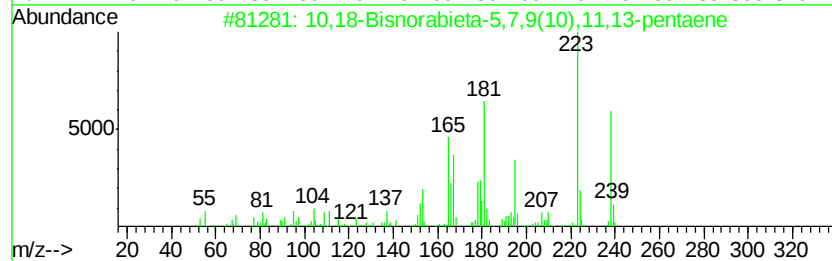
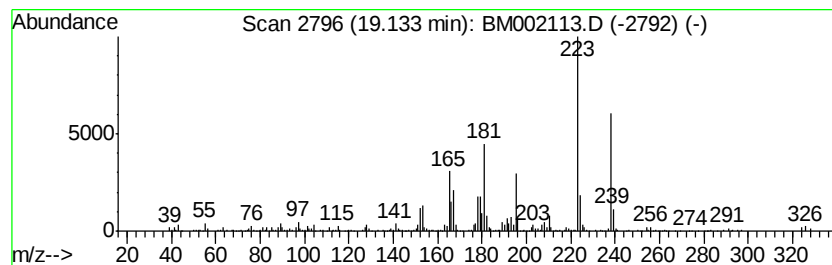
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 3

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.13 | 5.87 ng/ul | 444865 | Chrysene-d12     | 21.21 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 10,18-Bisnorabieta-5,7,9(10),11,... | 238 | C18H22   | 006566-19-4  | 99   |
| 2    |    |   | 12-Azabicyclo[9.2.2]pentadecan-1... | 223 | C14H25NO | 1000191-26-7 | 78   |
| 3    |    |   | Tricyclo[3.3.3.0(1,5)]undec-6-en... | 223 | C14H13N3 | 118894-71-6  | 78   |
| 4    |    |   | 3-(N,N-Diethylamino)carbazole       | 238 | C16H18N2 | 054994-31-9  | 58   |
| 5    |    |   | 2-Propenoic acid, 3-(3,4,5-trime... | 238 | C12H14O5 | 000090-50-6  | 53   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

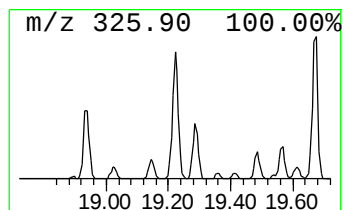
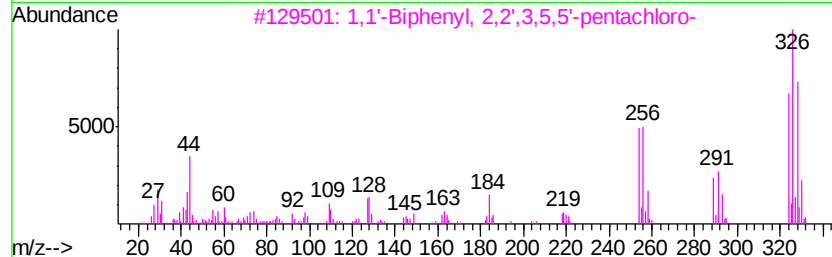
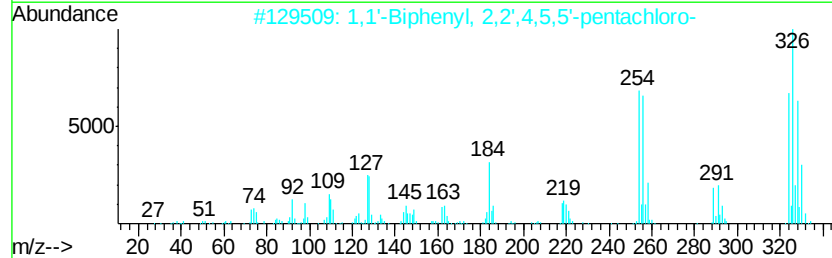
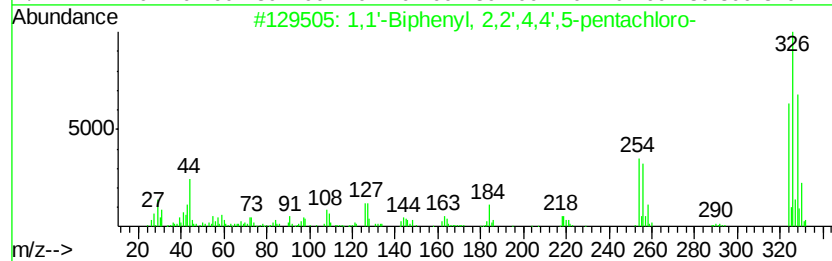
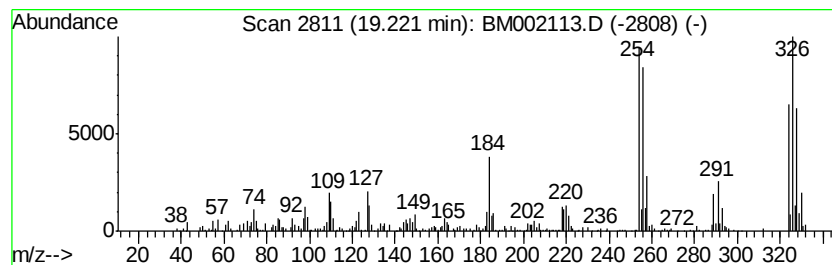
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 1,1'-Biphenyl, 2,2',4,4',5-... Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.22 | 4.56 ng/ul | 345673 | Chrysene-d12     | 21.21 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 99   |
| 2    |    |   | 1,1'-Biphenyl, 2,2',4,5,5'-penta... | 324 | C12H5Cl5 | 037680-73-2 | 99   |
| 3    |    |   | 1,1'-Biphenyl, 2,2',3,5,5'-penta... | 324 | C12H5Cl5 | 052663-61-3 | 98   |
| 4    |    |   | 1,1'-Biphenyl, 2,3',4,4',5-penta... | 324 | C12H5Cl5 | 031508-00-6 | 96   |
| 5    |    |   | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 96   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

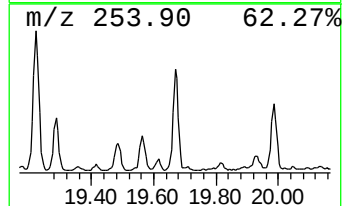
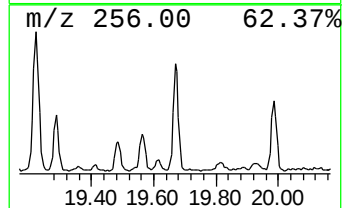
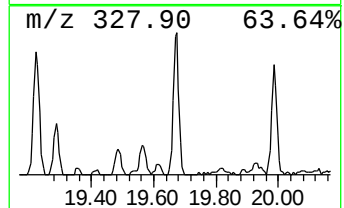
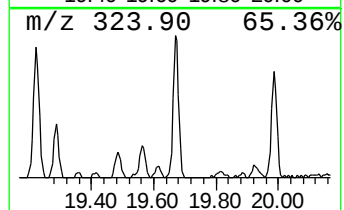
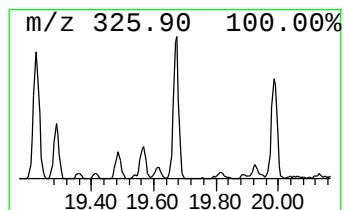
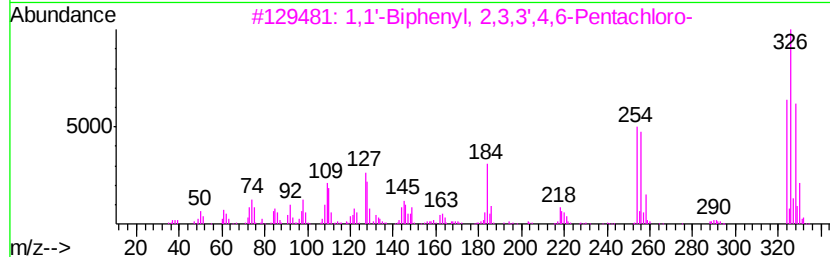
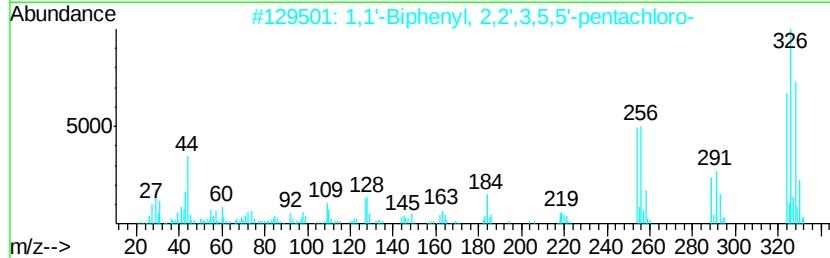
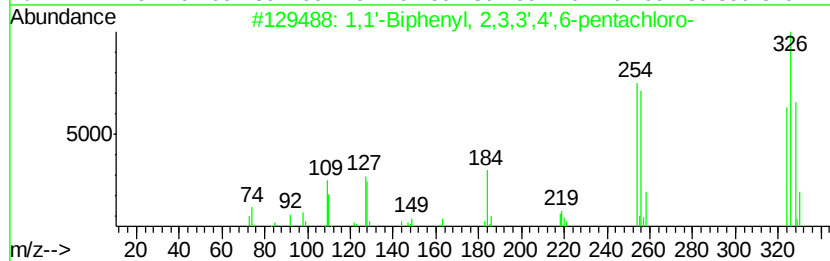
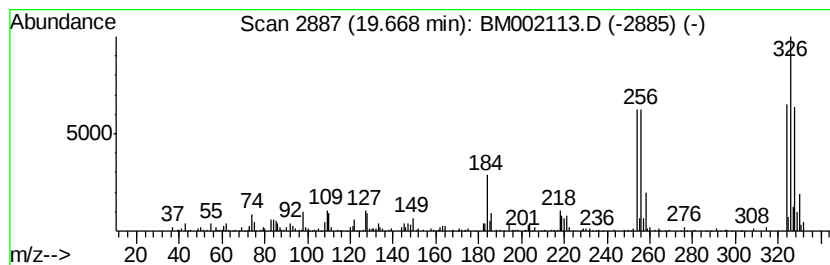
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 1,1'-Biphenyl, 2,3,3',4',6-... Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.67 | 4.14 ng/ul | 314117 | Chrysene-d12     | 21.21 |

| Hit# | of   | 5          | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
|------|------|------------|----------------------|-----|----------|-------------|------|
| 1    | 1,1' | -Biphenyl, | 2,3,3',4',6-penta... | 324 | C12H5Cl5 | 038380-03-9 | 99   |
| 2    | 1,1' | -Biphenyl, | 2,2',3,5,5'-penta... | 324 | C12H5Cl5 | 052663-61-3 | 96   |
| 3    | 1,1' | -Biphenyl, | 2,3,3',4,6-Pentac... | 324 | C12H5Cl5 | 074472-35-8 | 96   |
| 4    | 1,1' | -Biphenyl, | 2,2',4,4',6-Penta... | 324 | C12H5Cl5 | 039485-83-1 | 95   |
| 5    | 1,1' | -Biphenyl, | 2,3,3',4,4'-penta... | 324 | C12H5Cl5 | 032598-14-4 | 95   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

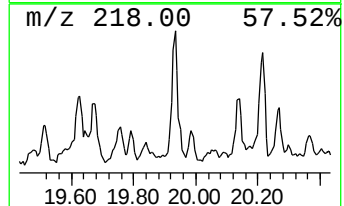
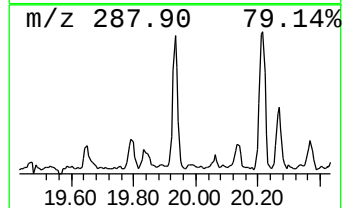
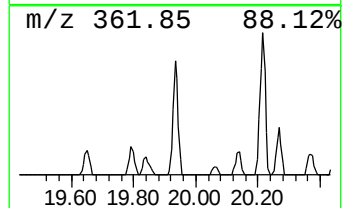
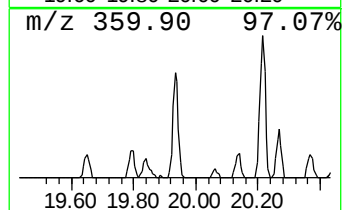
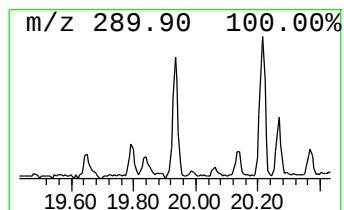
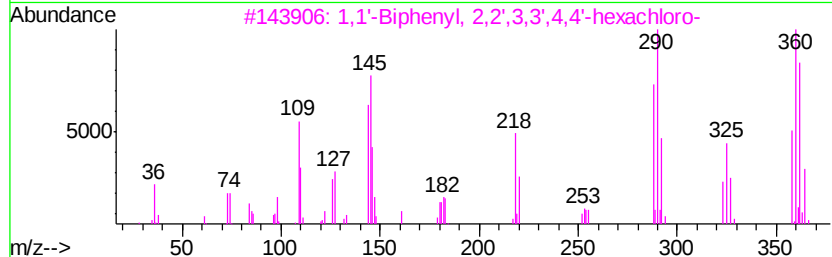
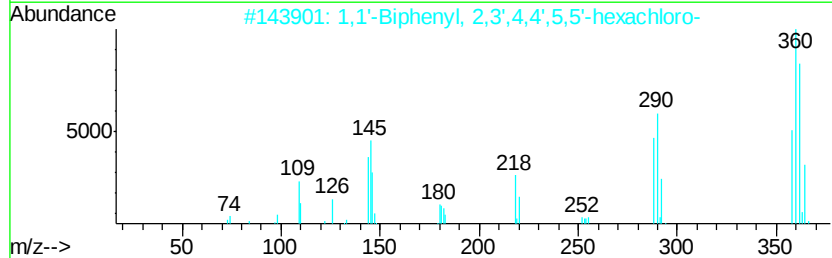
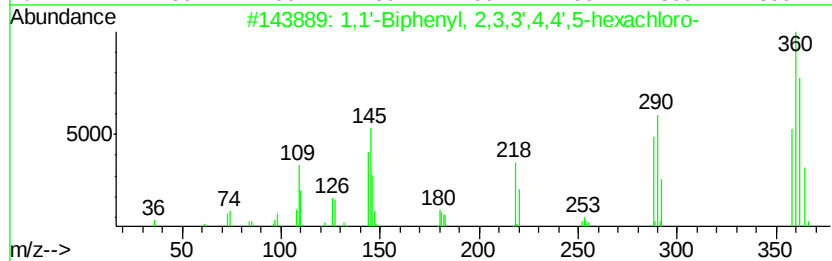
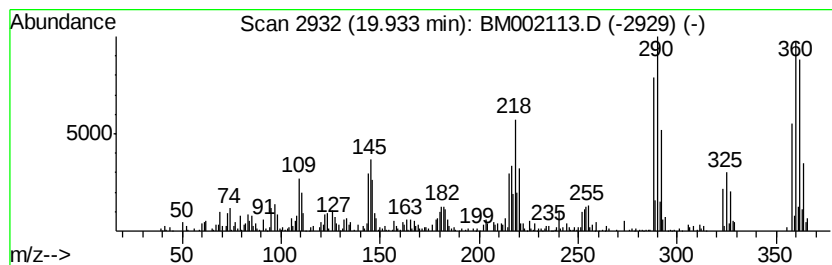
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 1,1'-Biphenyl, 2,3,3',4,4',... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.93 | 3.82 ng/ul | 289886 | Chrysene-d12     | 21.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2,3,3',4,4',5-hex... | 358 | C12H4Cl6 | 038380-08-4 | 99   |
| 2    |    | 1,1'-Biphenyl, 2,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 052663-72-6 | 97   |
| 3    |    | 1,1'-Biphenyl, 2,2',3,3',4,4'-he... | 358 | C12H4Cl6 | 038380-07-3 | 96   |
| 4    |    | 1,1'-Biphenyl, 2,2',3,4,5,6'-Hex... | 358 | C12H4Cl6 | 068194-15-0 | 96   |
| 5    |    | 1,1'-Biphenyl, 2,2',3,3',4,4'-he... | 358 | C12H4Cl6 | 038380-07-3 | 96   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

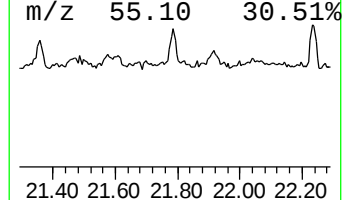
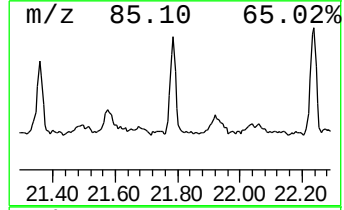
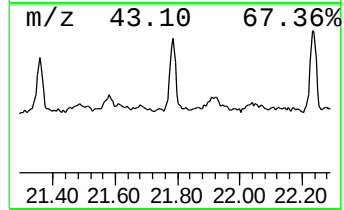
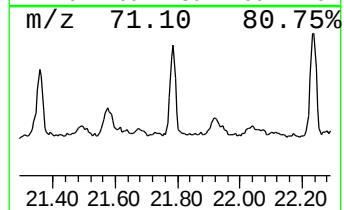
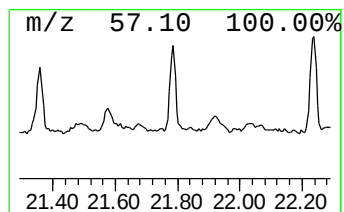
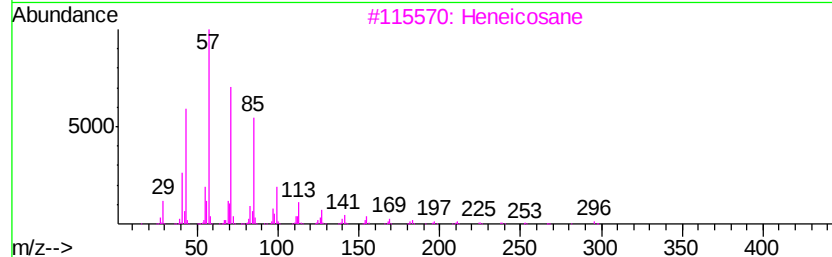
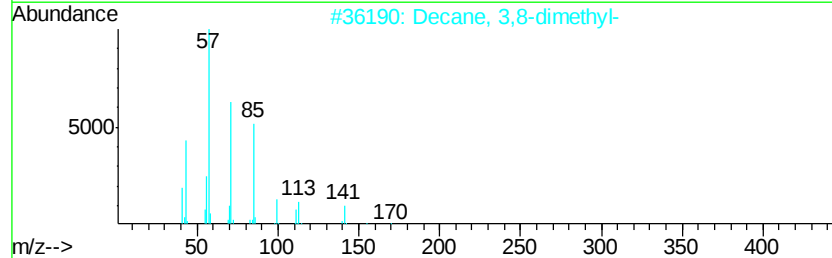
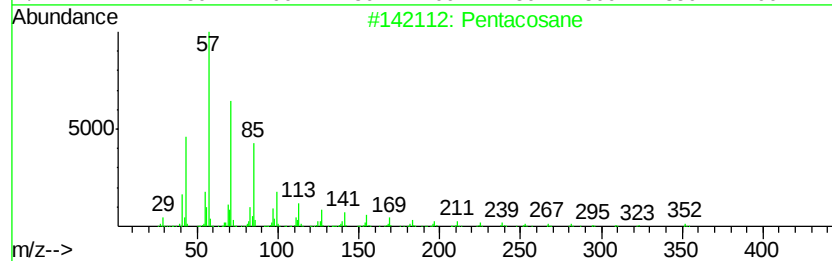
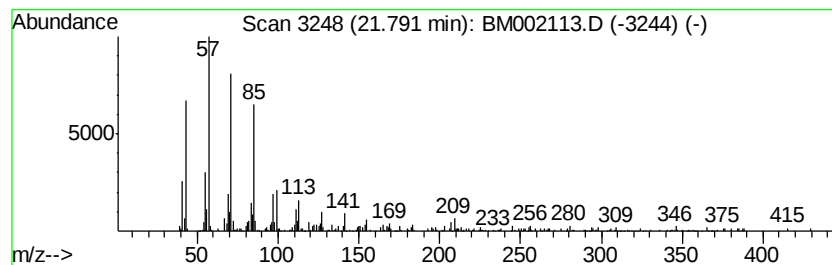
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.79 | 4.79 ng/ul | 363298 | Chrysene-d12     | 21.21 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Pentacosane           | 352 | C25H52  | 000629-99-2 | 90   |
| 2    |    | Decane, 3,8-dimethyl- | 170 | C12H26  | 017312-55-9 | 87   |
| 3    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 87   |
| 4    |    | Tetratriacontane      | 479 | C34H70  | 014167-59-0 | 87   |
| 5    |    | Nonadecane            | 268 | C19H40  | 000629-92-5 | 87   |



Data Path : Z:\HPCHEM1\BNA M\DATA\BM072815\  
Data File : BM002113.D  
Acq On : 29 Jul 2015 03:28  
Operator : TP/UM  
Sample : G3048-13  
Misc :  
ALS Vial : 21 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C01Y6

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM072715.M  
Quant Title : SVOA CALIBRATION

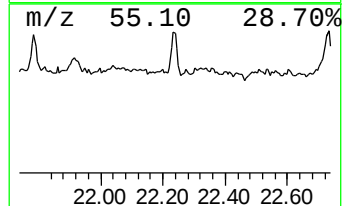
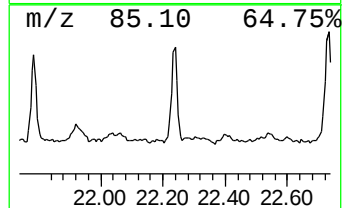
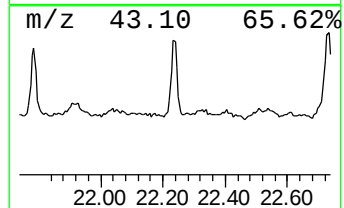
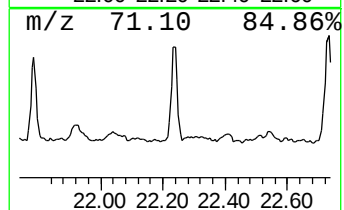
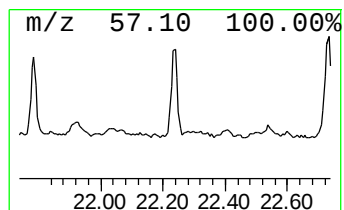
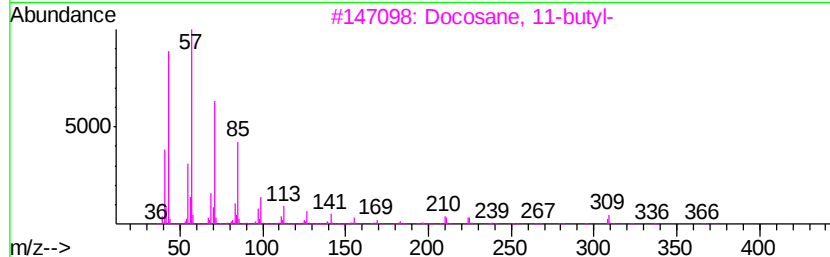
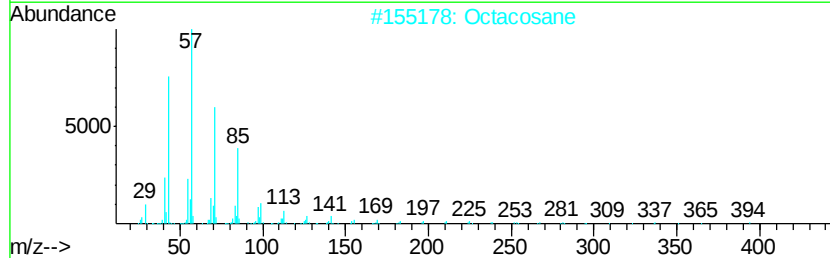
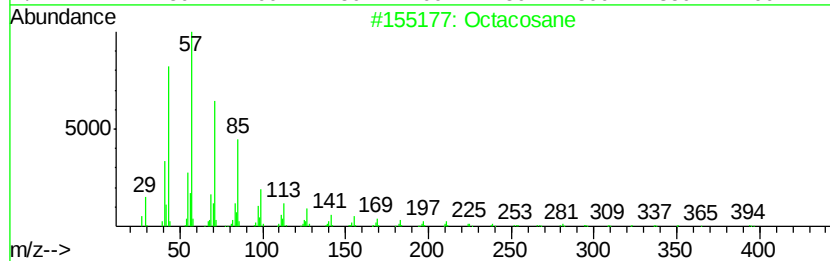
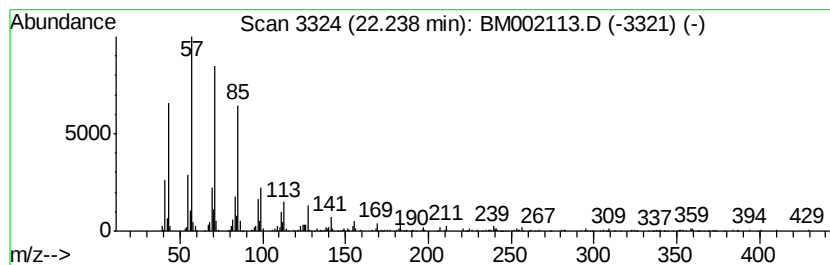
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.24 | 6.87 ng/ul | 521020 | Chrysene-d12     | 21.21 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Octacosane                       | 394 | C28H58  | 000630-02-4 | 94   |
| 2    |    |   | Octacosane                       | 394 | C28H58  | 000630-02-4 | 93   |
| 3    |    |   | Docosane, 11-butyl-              | 366 | C26H54  | 013475-76-8 | 93   |
| 4    |    |   | Heneicosane, 11-(1-ethylpropyl)- | 366 | C26H54  | 055282-11-6 | 91   |
| 5    |    |   | triacontane                      | 422 | C30H62  | 000638-68-6 | 91   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM072815\  
 Data File : BM002113.D  
 Acq On : 29 Jul 2015 03:28  
 Operator : TP/UM  
 Sample : G3048-13  
 Misc :  
 ALS Vial : 21 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C01Y6

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM072715.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| (DEL) Alkane: Str... | 11.40 | 2.0     | ng/ul | 53835    | 2                     | 10.39 | 534364  | 20.0 |
| unknown-01           | 14.08 | 2.8     | ng/ul | 101556   | 3                     | 14.26 | 715103  | 20.0 |
| (DEL) Alkane: Str... | 15.52 | 4.0     | ng/ul | 142234   | 3                     | 14.26 | 715103  | 20.0 |
| (DEL) Alkane: Str... | 16.00 | 10.3    | ng/ul | 454713   | 4                     | 17.01 | 880803  | 20.0 |
| (DEL) Alkane: Str... | 16.82 | 4.9     | ng/ul | 217002   | 4                     | 17.01 | 880803  | 20.0 |
| (DEL) Alkane: Str... | 17.42 | 2.2     | ng/ul | 96647    | 4                     | 17.01 | 880803  | 20.0 |
| 1,1'-Biphenyl, 2,... | 18.08 | 3.9     | ng/ul | 172906   | 4                     | 17.01 | 880803  | 20.0 |
| 4b,8-Dimethyl-2-i... | 18.63 | 3.5     | ng/ul | 155076   | 4                     | 17.01 | 880803  | 20.0 |
| 10,18-Bisnorabiet... | 19.13 | 5.9     | ng/ul | 444865   | 5                     | 21.21 | 1516270 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.22 | 4.6     | ng/ul | 345673   | 5                     | 21.21 | 1516270 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.67 | 4.1     | ng/ul | 314117   | 5                     | 21.21 | 1516270 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.93 | 3.8     | ng/ul | 289886   | 5                     | 21.21 | 1516270 | 20.0 |
| (DEL) Alkane: Str... | 21.79 | 4.8     | ng/ul | 363298   | 5                     | 21.21 | 1516270 | 20.0 |
| (DEL) Alkane: Str... | 22.24 | 6.9     | ng/ul | 521020   | 5                     | 21.21 | 1516270 | 20.0 |