

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
 Data File : BG018214.D  
 Acq On : 1 Aug 2015 11:03  
 Operator : TP/UM  
 Sample : G3073-09  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C01Q2

## 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\_G\METHODS\SOM02.2-EPA-BG073115.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         | 3.313       | 102           | 106         | 113          | rVB2     | 41023          | 53369         | 1.67%           | 0.136%        |
| 2         | 6.621       | 665           | 669         | 675          | rBV5     | 16256          | 34114         | 1.07%           | 0.087%        |
| 3         | 6.686       | 675           | 680         | 690          | rVB      | 122852         | 205493        | 6.43%           | 0.524%        |
| 4         | 6.827       | 696           | 704         | 710          | rBV      | 90956          | 149660        | 4.68%           | 0.382%        |
| 5         | 7.027       | 733           | 738         | 744          | rVB      | 118525         | 191259        | 5.98%           | 0.488%        |
| 6         | 7.138       | 749           | 757         | 760          | rBV3     | 24298          | 44116         | 1.38%           | 0.113%        |
| 7         | 7.403       | 793           | 802         | 806          | rVV10    | 20169          | 49694         | 1.55%           | 0.127%        |
| 8         | 7.485       | 806           | 816         | 823          | rVB2     | 230203         | 398553        | 12.46%          | 1.017%        |
| 9         | 7.673       | 843           | 848         | 851          | rVB5     | 25667          | 35458         | 1.11%           | 0.090%        |
| 10        | 7.749       | 857           | 861         | 871          | rBV9     | 28958          | 66711         | 2.09%           | 0.170%        |
| 11        | 7.978       | 896           | 900         | 904          | rVV5     | 17522          | 34541         | 1.08%           | 0.088%        |
| 12        | 8.031       | 905           | 909         | 914          | rVB5     | 41265          | 66357         | 2.07%           | 0.169%        |
| 13        | 8.125       | 919           | 925         | 930          | rBV5     | 32026          | 59427         | 1.86%           | 0.152%        |
| 14        | 8.213       | 935           | 940         | 947          | rVB      | 141528         | 210612        | 6.59%           | 0.537%        |
| 15        | 8.302       | 947           | 955         | 959          | rBV5     | 57504          | 112584        | 3.52%           | 0.287%        |
| 16        | 8.584       | 992           | 1003        | 1008         | rBV8     | 76650          | 192711        | 6.03%           | 0.492%        |
| 17        | 8.642       | 1008          | 1013        | 1017         | rVV      | 116170         | 186023        | 5.82%           | 0.475%        |
| 18        | 8.801       | 1034          | 1040        | 1041         | rBV3     | 51580          | 86032         | 2.69%           | 0.220%        |
| 19        | 8.830       | 1041          | 1045        | 1051         | rVB5     | 75503          | 158518        | 4.96%           | 0.405%        |
| 20        | 8.936       | 1051          | 1063        | 1071         | rBV5     | 41450          | 170899        | 5.34%           | 0.436%        |
| 21        | 9.124       | 1089          | 1095        | 1100         | rBV6     | 77868          | 155388        | 4.86%           | 0.397%        |
| 22        | 9.189       | 1100          | 1106        | 1113         | rVB5     | 101454         | 196404        | 6.14%           | 0.501%        |
| 23        | 9.359       | 1128          | 1135        | 1140         | rBV4     | 109999         | 233585        | 7.30%           | 0.596%        |
| 24        | 9.400       | 1140          | 1142        | 1145         | rVV      | 45022          | 48732         | 1.52%           | 0.124%        |
| 25        | 9.447       | 1145          | 1150        | 1159         | rVB7     | 103414         | 230631        | 7.21%           | 0.589%        |
| 26        | 9.677       | 1184          | 1189        | 1193         | rBV7     | 38503          | 71314         | 2.23%           | 0.182%        |
| 27        | 9.864       | 1214          | 1221        | 1224         | rBV7     | 51501          | 96915         | 3.03%           | 0.247%        |
| 28        | 9.906       | 1224          | 1228        | 1235         | rVB      | 164920         | 249085        | 7.79%           | 0.636%        |
| 29        | 9.970       | 1235          | 1239        | 1243         | rBV3     | 46280          | 85661         | 2.68%           | 0.219%        |
| 30        | 10.135      | 1263          | 1267        | 1270         | rBV5     | 60269          | 94770         | 2.96%           | 0.242%        |
| 31        | 10.176      | 1271          | 1274        | 1278         | rVV      | 40904          | 59149         | 1.85%           | 0.151%        |
| 32        | 10.270      | 1278          | 1290        | 1295         | rVV      | 369516         | 769289        | 24.05%          | 1.963%        |
| 33        | 10.323      | 1295          | 1299        | 1306         | rVV2     | 96203          | 216771        | 6.78%           | 0.553%        |
| 34        | 10.393      | 1306          | 1311        | 1313         | rVV3     | 78804          | 139521        | 4.36%           | 0.356%        |

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Integrator: RTE  
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 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\_G\METHODS\SOM02.2-EPA-BG073115.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 10.446 | 1316 | 1320 | 1324 | rVV2 | 181193 | 284911  | 8.91%  | 0.727% |
| 36 | 10.499 | 1326 | 1329 | 1335 | rVB7 | 31823  | 58655   | 1.83%  | 0.150% |
| 37 | 10.675 | 1353 | 1359 | 1363 | rBV6 | 58770  | 114975  | 3.60%  | 0.293% |
| 38 | 10.763 | 1370 | 1374 | 1378 | rVB3 | 67223  | 98004   | 3.06%  | 0.250% |
| 39 | 10.805 | 1378 | 1381 | 1384 | rBV5 | 34725  | 46575   | 1.46%  | 0.119% |
| 40 | 10.957 | 1402 | 1407 | 1415 | rVB7 | 125538 | 338196  | 10.57% | 0.863% |
| 41 | 11.310 | 1462 | 1467 | 1471 | rBV3 | 320043 | 536213  | 16.77% | 1.368% |
| 42 | 11.563 | 1506 | 1510 | 1515 | rVB5 | 81873  | 120979  | 3.78%  | 0.309% |
| 43 | 12.021 | 1583 | 1588 | 1591 | rBV3 | 99811  | 182775  | 5.72%  | 0.466% |
| 44 | 12.091 | 1597 | 1600 | 1602 | rBV4 | 35242  | 44558   | 1.39%  | 0.114% |
| 45 | 12.432 | 1653 | 1658 | 1665 | rVB6 | 97809  | 199506  | 6.24%  | 0.509% |
| 46 | 12.497 | 1665 | 1669 | 1673 | rBV7 | 85276  | 133177  | 4.16%  | 0.340% |
| 47 | 12.573 | 1680 | 1682 | 1688 | rVB5 | 129885 | 182805  | 5.72%  | 0.467% |
| 48 | 12.649 | 1690 | 1695 | 1698 | rVV6 | 70068  | 143209  | 4.48%  | 0.365% |
| 49 | 12.696 | 1699 | 1703 | 1707 | rVV  | 390679 | 530283  | 16.58% | 1.353% |
| 50 | 12.743 | 1708 | 1711 | 1717 | rVB5 | 98499  | 160195  | 5.01%  | 0.409% |
| 51 | 12.926 | 1738 | 1742 | 1745 | rBV  | 153587 | 214702  | 6.71%  | 0.548% |
| 52 | 13.078 | 1765 | 1768 | 1774 | rVB6 | 94882  | 124732  | 3.90%  | 0.318% |
| 53 | 13.578 | 1848 | 1853 | 1857 | rVB2 | 466804 | 685524  | 21.44% | 1.749% |
| 54 | 13.625 | 1858 | 1861 | 1863 | rVV  | 132550 | 158284  | 4.95%  | 0.404% |
| 55 | 13.660 | 1863 | 1867 | 1871 | rVB2 | 427232 | 582528  | 18.21% | 1.487% |
| 56 | 13.854 | 1896 | 1900 | 1904 | rBV  | 158464 | 236799  | 7.40%  | 0.604% |
| 57 | 13.989 | 1919 | 1923 | 1927 | rVB2 | 246373 | 300089  | 9.38%  | 0.766% |
| 58 | 14.077 | 1933 | 1938 | 1942 | rBV4 | 150856 | 253259  | 7.92%  | 0.646% |
| 59 | 14.124 | 1942 | 1946 | 1948 | rVV3 | 144638 | 208895  | 6.53%  | 0.533% |
| 60 | 14.165 | 1949 | 1953 | 1961 | rVB2 | 400063 | 698374  | 21.84% | 1.782% |
| 61 | 14.706 | 2041 | 2045 | 2050 | rVB4 | 232827 | 324420  | 10.14% | 0.828% |
| 62 | 14.953 | 2083 | 2087 | 2092 | rVB3 | 203821 | 282509  | 8.83%  | 0.721% |
| 63 | 15.041 | 2099 | 2102 | 2105 | rVB  | 149449 | 168341  | 5.26%  | 0.430% |
| 64 | 15.164 | 2120 | 2123 | 2128 | rVB  | 297740 | 396847  | 12.41% | 1.013% |
| 65 | 15.217 | 2129 | 2132 | 2136 | rBV4 | 91444  | 136587  | 4.27%  | 0.349% |
| 66 | 15.446 | 2167 | 2171 | 2176 | rVB  | 323159 | 464960  | 14.54% | 1.187% |
| 67 | 15.652 | 2203 | 2206 | 2214 | rVB5 | 104998 | 154240  | 4.82%  | 0.394% |
| 68 | 15.928 | 2245 | 2253 | 2258 | rBV  | 517208 | 1083285 | 33.87% | 2.765% |
| 69 | 16.698 | 2381 | 2384 | 2388 | rBV3 | 89371  | 107870  | 3.37%  | 0.275% |
| 70 | 16.745 | 2388 | 2392 | 2400 | rVB2 | 294328 | 488722  | 15.28% | 1.247% |
| 71 | 16.921 | 2418 | 2422 | 2426 | rBV2 | 364719 | 506624  | 15.84% | 1.293% |

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BNA\_G  
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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\_G\METHODS\SOM02.2-EPA-BG073115.M  
Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 16.962 | 2426 | 2429 | 2434 | rVB  | 155043  | 207346  | 6.48%   | 0.529% |
| 73  | 17.021 | 2435 | 2439 | 2443 | rVB  | 234473  | 313248  | 9.79%   | 0.799% |
| 74  | 17.432 | 2507 | 2509 | 2514 | rVB3 | 103944  | 116102  | 3.63%   | 0.296% |
| 75  | 17.838 | 2575 | 2578 | 2583 | rVB3 | 89198   | 126470  | 3.95%   | 0.323% |
| 76  | 18.554 | 2696 | 2700 | 2704 | rVB  | 245144  | 310829  | 9.72%   | 0.793% |
| 77  | 18.772 | 2735 | 2737 | 2741 | rBV4 | 40027   | 43371   | 1.36%   | 0.111% |
| 78  | 18.995 | 2770 | 2775 | 2781 | rBV  | 359622  | 492080  | 15.39%  | 1.256% |
| 79  | 19.054 | 2781 | 2785 | 2791 | rVB2 | 305024  | 394643  | 12.34%  | 1.007% |
| 80  | 19.142 | 2796 | 2800 | 2805 | rVB5 | 90962   | 134873  | 4.22%   | 0.344% |
| 81  | 19.330 | 2828 | 2832 | 2834 | rBV2 | 288625  | 381305  | 11.92%  | 0.973% |
| 82  | 19.359 | 2834 | 2837 | 2841 | rVB  | 331169  | 402986  | 12.60%  | 1.028% |
| 83  | 19.594 | 2874 | 2877 | 2883 | rVB3 | 81372   | 118471  | 3.70%   | 0.302% |
| 84  | 19.859 | 2919 | 2922 | 2925 | rVB2 | 83615   | 93929   | 2.94%   | 0.240% |
| 85  | 20.135 | 2967 | 2969 | 2972 | rBV3 | 53136   | 48585   | 1.52%   | 0.124% |
| 86  | 20.393 | 3010 | 3013 | 3016 | rVB  | 149514  | 150760  | 4.71%   | 0.385% |
| 87  | 20.858 | 3088 | 3092 | 3097 | rVB2 | 416119  | 478938  | 14.98%  | 1.222% |
| 88  | 21.128 | 3132 | 3138 | 3142 | rBV2 | 512196  | 883175  | 27.62%  | 2.254% |
| 89  | 21.292 | 3162 | 3166 | 3170 | rBV  | 819528  | 880866  | 27.54%  | 2.248% |
| 90  | 21.633 | 3220 | 3224 | 3227 | rBV  | 456733  | 543706  | 17.00%  | 1.388% |
| 91  | 21.715 | 3234 | 3238 | 3242 | rVB  | 1287917 | 1427765 | 44.64%  | 3.644% |
| 92  | 21.774 | 3244 | 3248 | 3253 | rBV2 | 780593  | 1002631 | 31.35%  | 2.559% |
| 93  | 21.927 | 3270 | 3274 | 3277 | rBV3 | 440092  | 580428  | 18.15%  | 1.481% |
| 94  | 22.168 | 3311 | 3315 | 3319 | rVB  | 1502042 | 1873597 | 58.58%  | 4.781% |
| 95  | 22.661 | 3394 | 3399 | 3410 | rVB  | 2163576 | 3198146 | 100.00% | 8.162% |
| 96  | 23.219 | 3488 | 3494 | 3500 | rVB  | 1647313 | 2704896 | 84.58%  | 6.903% |
| 97  | 23.313 | 3507 | 3510 | 3516 | rVB  | 261401  | 434379  | 13.58%  | 1.109% |
| 98  | 23.848 | 3594 | 3601 | 3609 | rVB  | 1431483 | 2680014 | 83.80%  | 6.839% |
| 99  | 24.577 | 3718 | 3725 | 3733 | rVB2 | 839871  | 1828316 | 57.17%  | 4.666% |
| 100 | 25.435 | 3864 | 3871 | 3880 | rVB  | 592575  | 1427339 | 44.63%  | 3.643% |

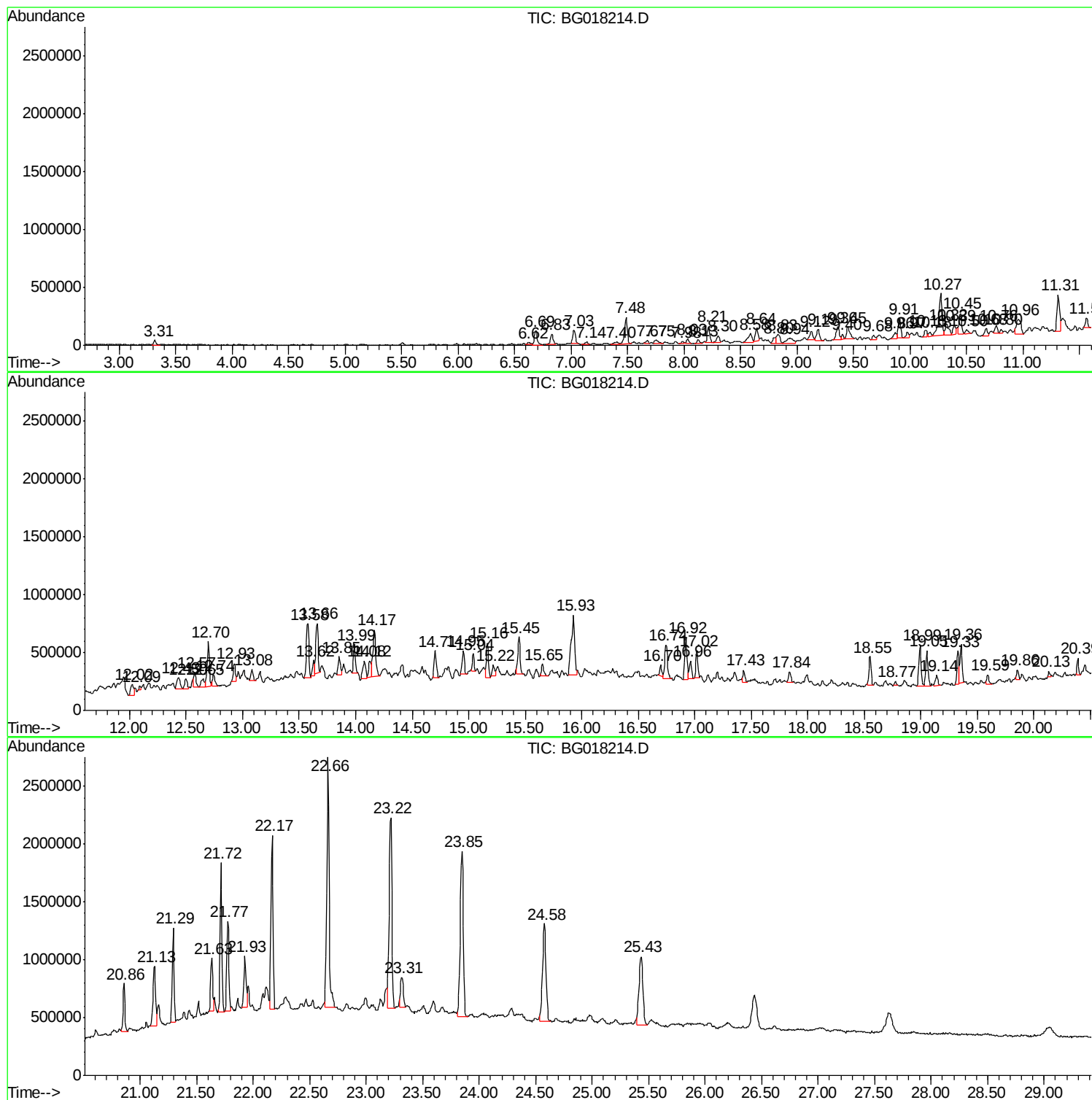
Sum of corrected areas: 39185147

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Instrument :  
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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

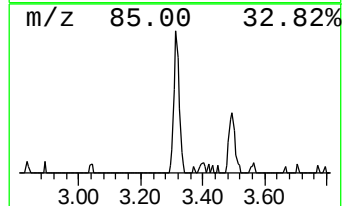
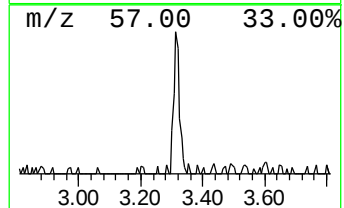
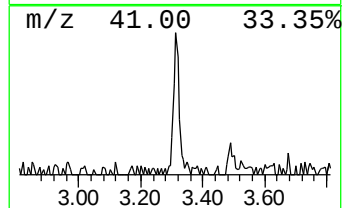
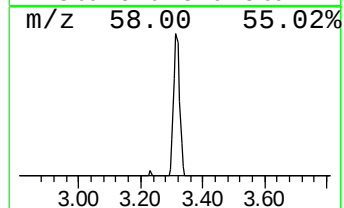
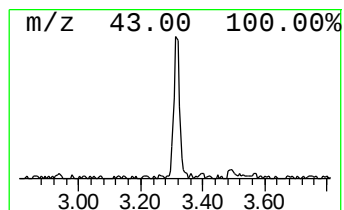
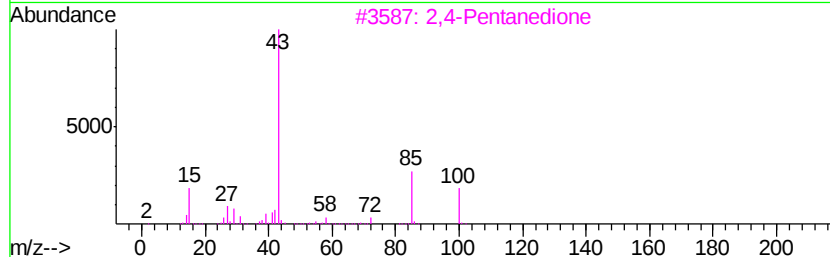
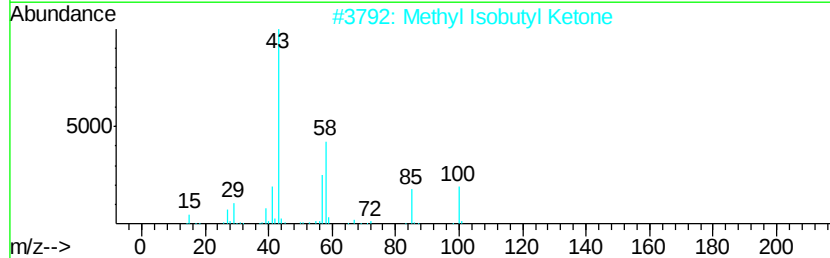
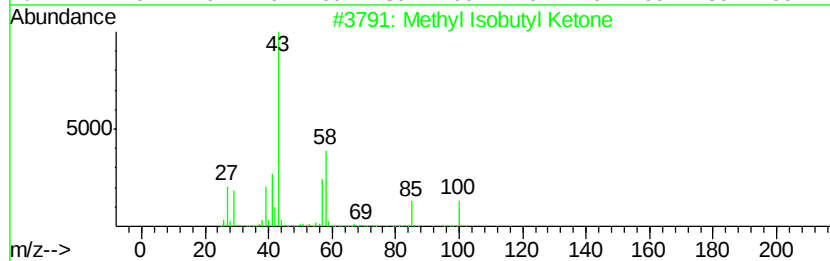
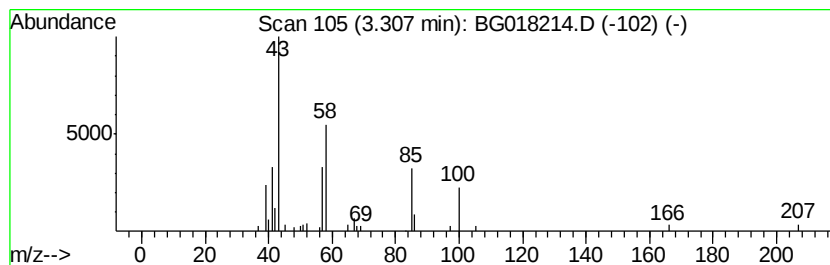
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 Methyl Isobutyl Ketone Concentration Rank 52

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 3.31 | 2.68 ng/ul | 53369 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Methyl Isobutyl Ketone | 100 | C6H12O  | 000108-10-1 | 56   |
| 2    |    | Methyl Isobutyl Ketone | 100 | C6H12O  | 000108-10-1 | 53   |
| 3    |    | 2,4-Pentanedione       | 100 | C5H8O2  | 000123-54-6 | 47   |
| 4    |    | 2,4-Pentanedione       | 100 | C5H8O2  | 000123-54-6 | 38   |
| 5    |    | 2,4-Pentanedione       | 100 | C5H8O2  | 000123-54-6 | 38   |



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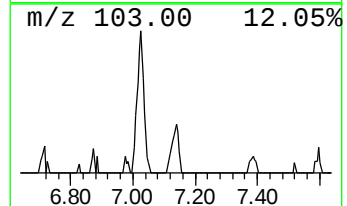
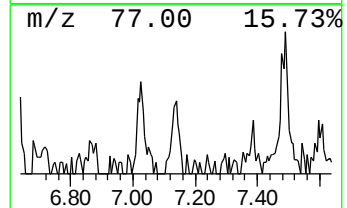
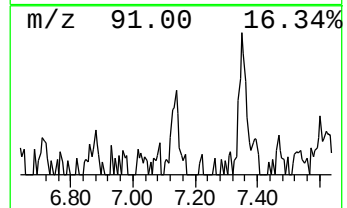
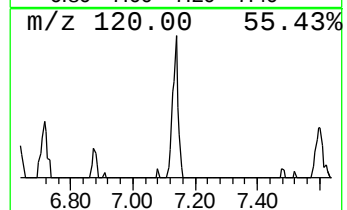
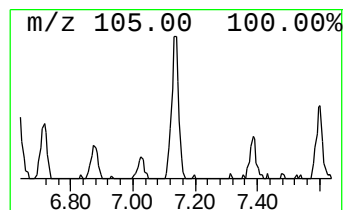
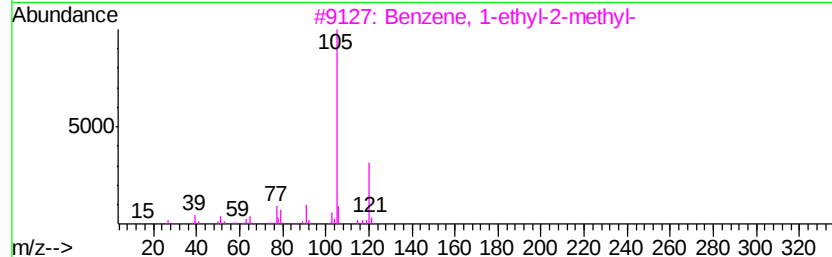
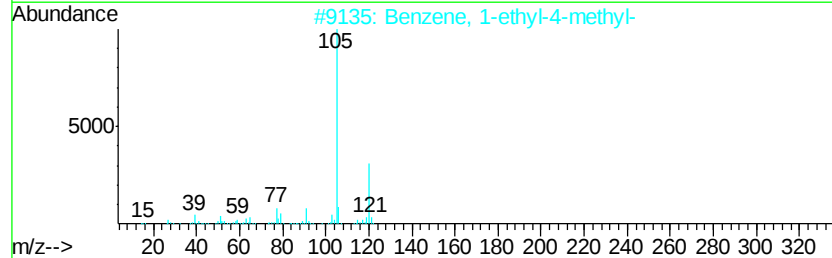
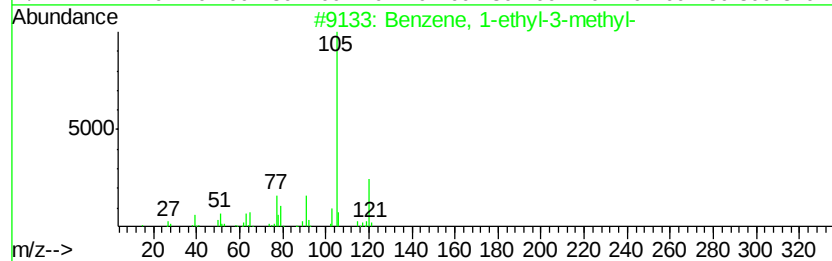
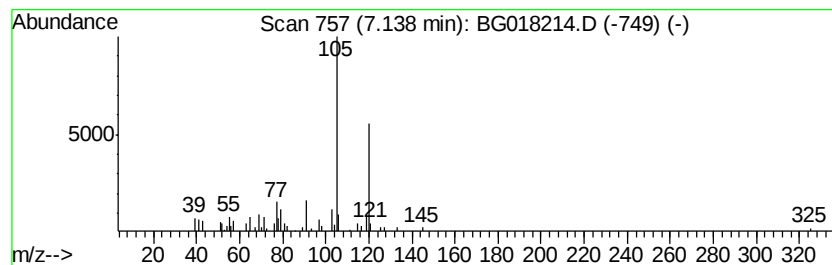
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 57

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 7.14 | 2.21 ng/ul | 44116 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 87   |
| 2    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 87   |
| 3    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 83   |
| 4    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 83   |
| 5    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 81   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

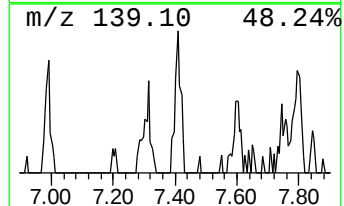
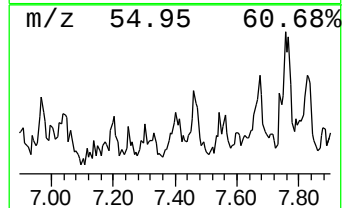
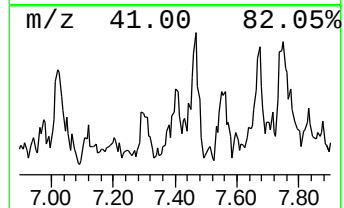
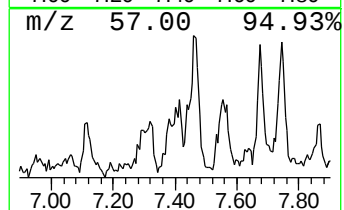
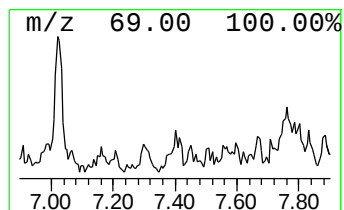
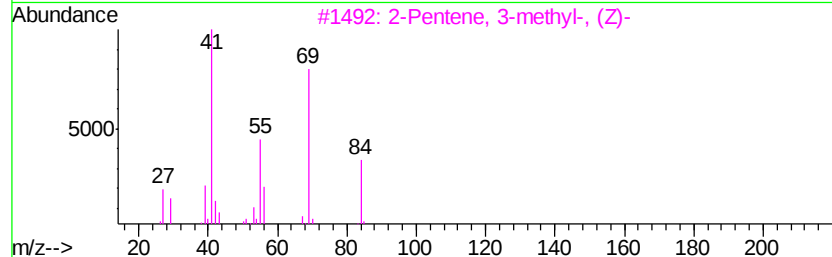
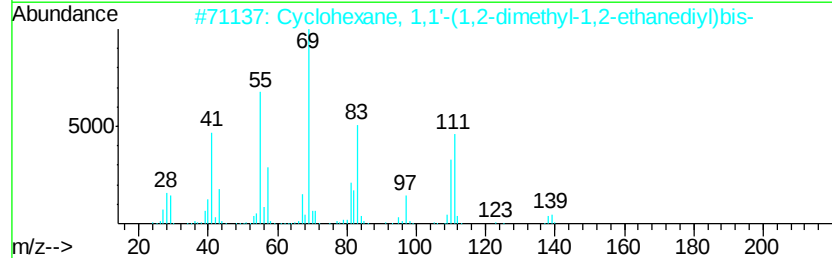
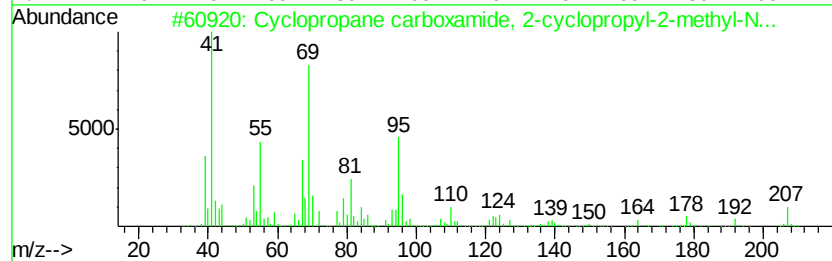
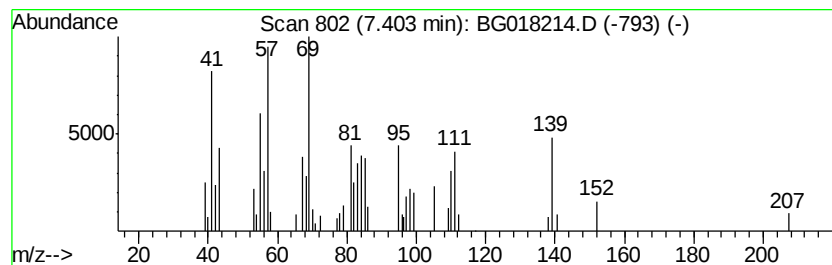
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 unknown-01 Concentration Rank 54

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 7.40 | 2.49 ng/ul | 49694 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Cyclopropane carboxamide, 2-cycl... | 207 | C13H21NO | 331416-19-4 | 38   |
| 2    |    | Cyclohexane, 1,1'-(1,2-dimethyl-... | 222 | C16H30   | 074663-71-1 | 35   |
| 3    |    | 2-Pentene, 3-methyl-, (Z)-          | 84  | C6H12    | 000922-62-3 | 25   |
| 4    |    | 2-Pentene, 3-methyl-, (Z)-          | 84  | C6H12    | 000922-62-3 | 25   |
| 5    |    | Pentane, 3-methylene-               | 84  | C6H12    | 000760-21-4 | 25   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleID :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

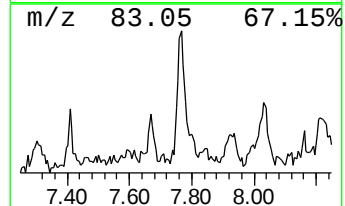
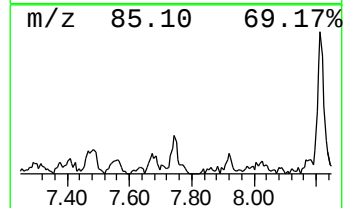
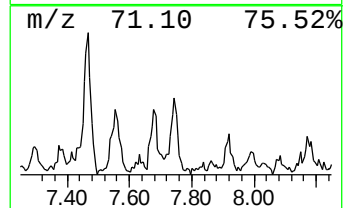
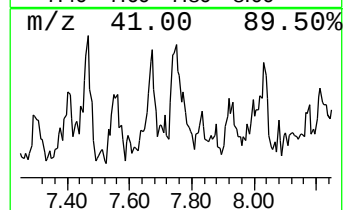
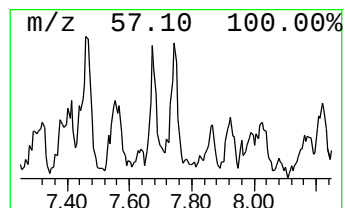
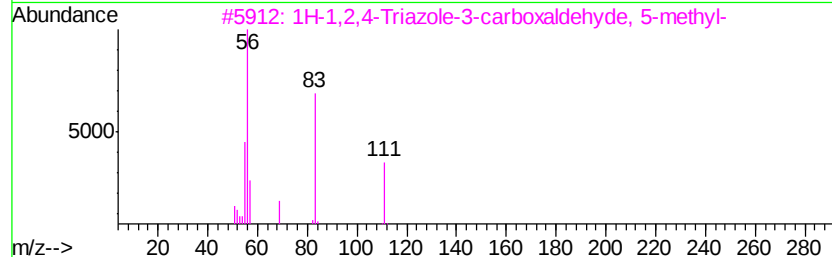
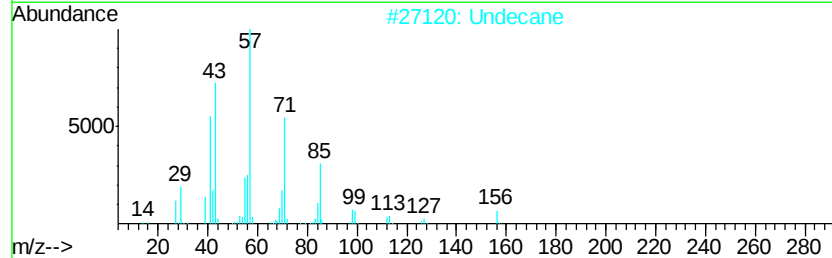
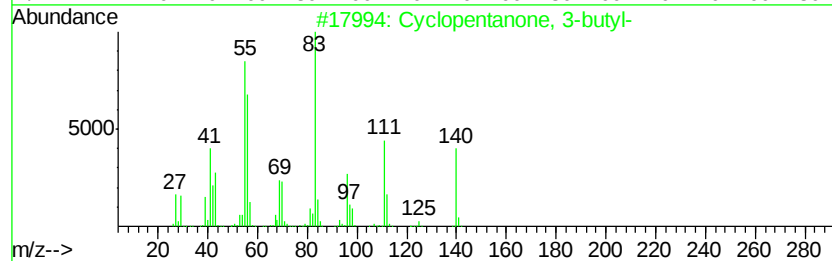
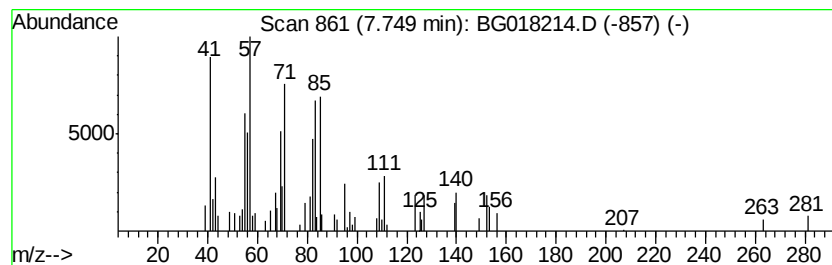
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 unknown-02 Concentration Rank 45

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 7.75 | 3.35 ng/ul | 66711 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopentanone, 3-butyl-            | 140 | C9H16O  | 057283-81-5 | 30   |
| 2    |    | Undecane                            | 156 | C11H24  | 001120-21-4 | 27   |
| 3    |    | 1H-1,2,4-Triazole-3-carboxaldehy... | 111 | C4H5N3O | 056804-98-9 | 18   |
| 4    |    | 1,3-Cyclohexanedione, 4,4-dimethyl- | 140 | C8H12O2 | 000562-46-9 | 14   |
| 5    |    | 2-Hexene, 1-(pentyloxy)-, (E)-      | 170 | C11H22O | 056052-82-5 | 14   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleID :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

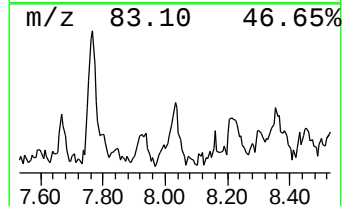
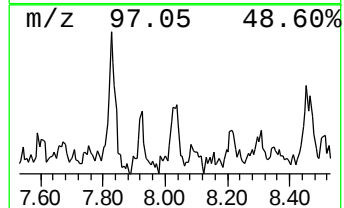
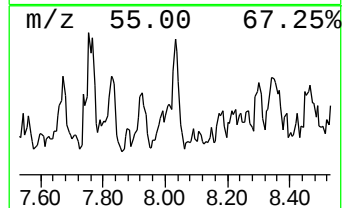
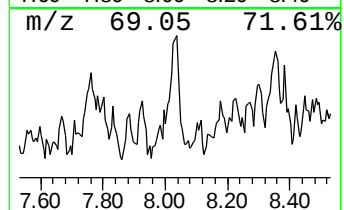
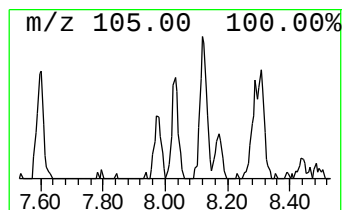
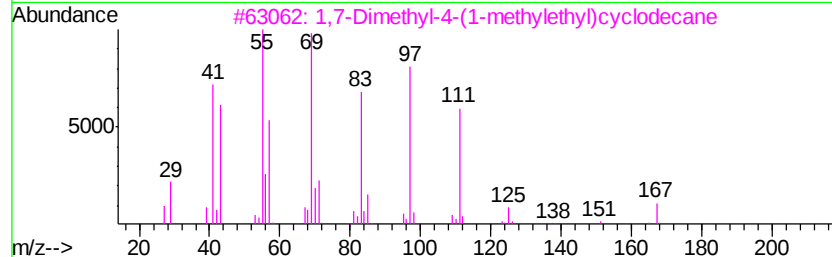
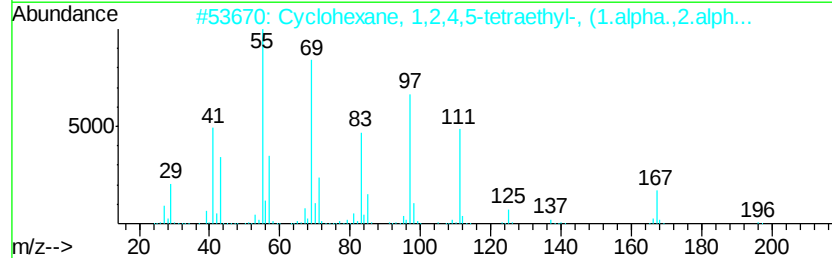
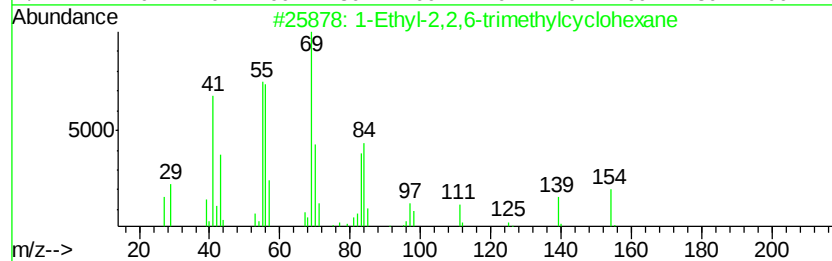
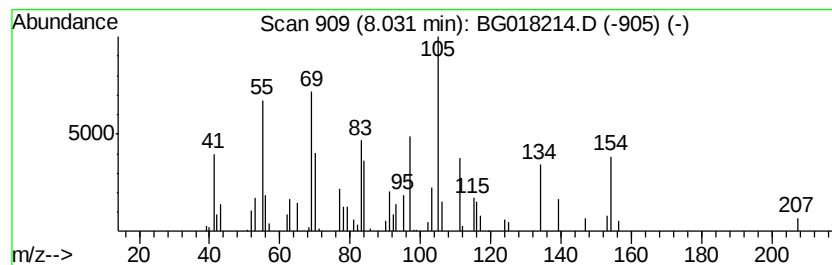
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Cyclic8.03 Concentration Rank 46

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 8.03 | 3.33 ng/ul | 66357 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1-Ethyl-2,2,6-trimethylcyclohexane  | 154 | C11H22  | 071186-27-1 | 28   |
| 2    |    | Cyclohexane, 1,2,4,5-tetraethyl-... | 196 | C14H28  | 061142-24-3 | 27   |
| 3    |    | 1,7-Dimethyl-4-(1-methylethyl)cy... | 210 | C15H30  | 000645-10-3 | 25   |
| 4    |    | 1-Dodecene                          | 168 | C12H24  | 000112-41-4 | 25   |
| 5    |    | 1,7-Dimethyl-4-(1-methylethyl)cy... | 210 | C15H30  | 000645-10-3 | 25   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
C01Q2

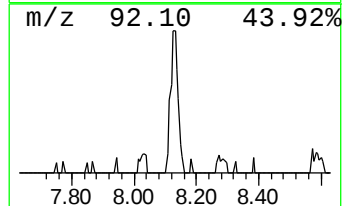
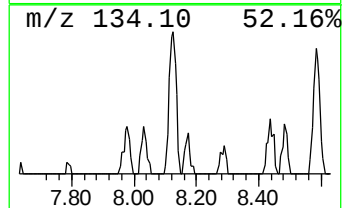
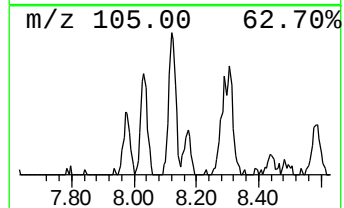
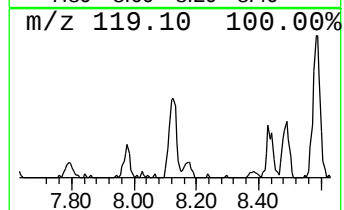
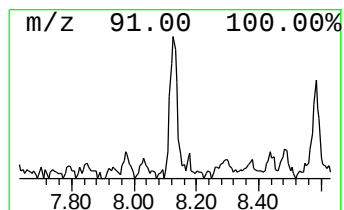
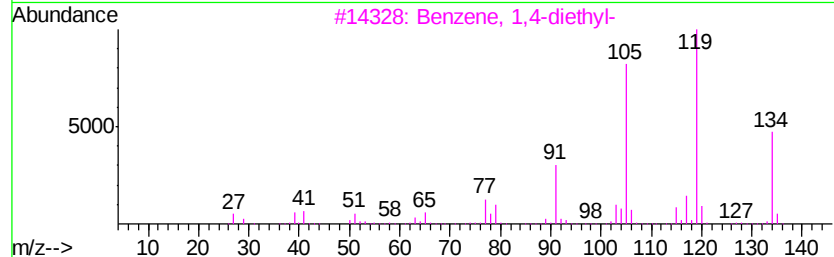
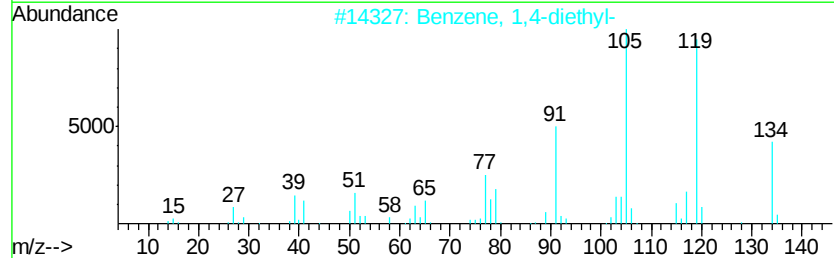
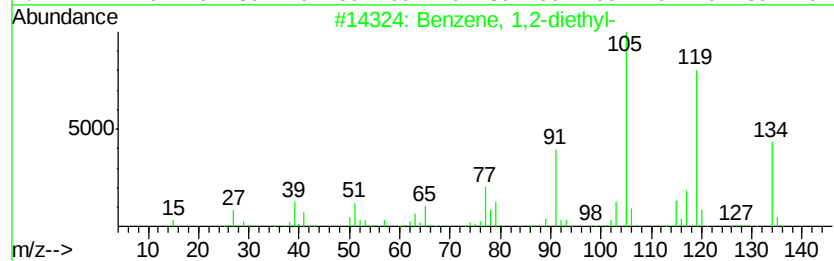
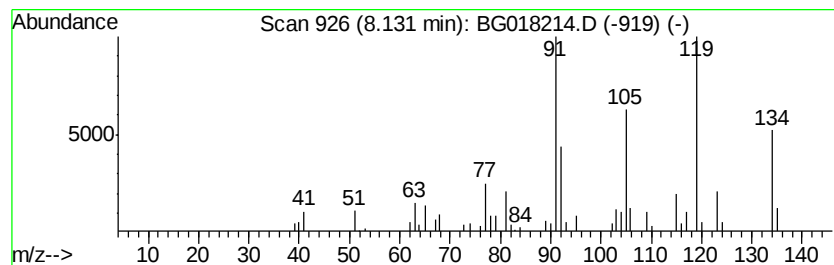
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 Benzene, 1,2-diethyl- Concentration Rank 50

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 8.13 | 2.98 ng/ul | 59427 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2-diethyl-               | 134 | C10H14  | 000135-01-3 | 78   |
| 2    |    | Benzene, 1,4-diethyl-               | 134 | C10H14  | 000105-05-5 | 76   |
| 3    |    | Benzene, 1,4-diethyl-               | 134 | C10H14  | 000105-05-5 | 70   |
| 4    |    | 1,3-Cyclopentadiene, 1,2,3,4-tet... | 134 | C10H14  | 076089-59-3 | 70   |
| 5    |    | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 70   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

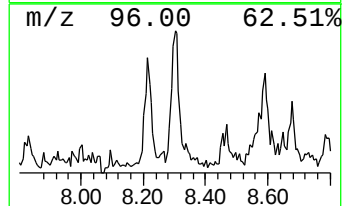
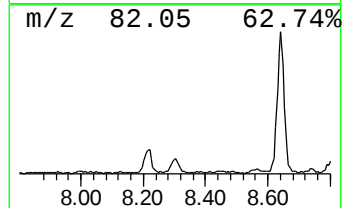
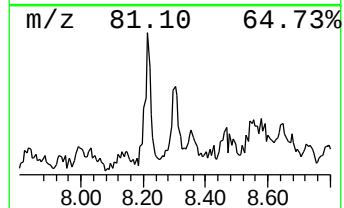
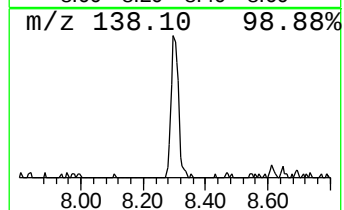
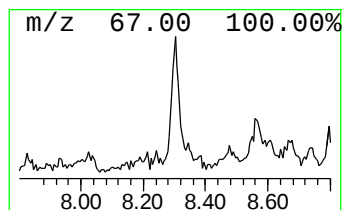
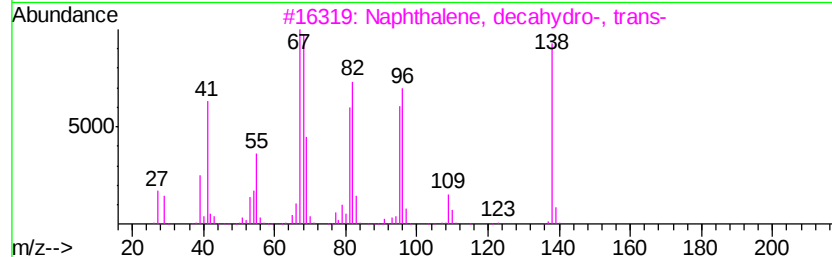
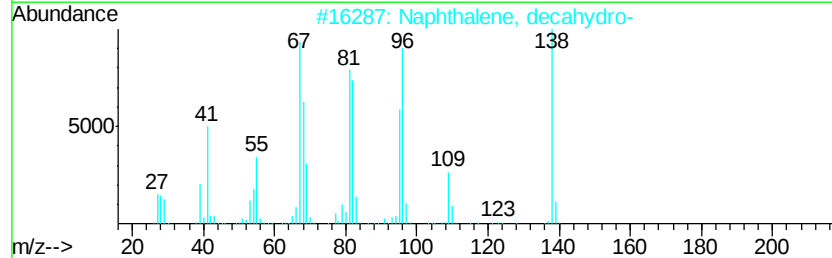
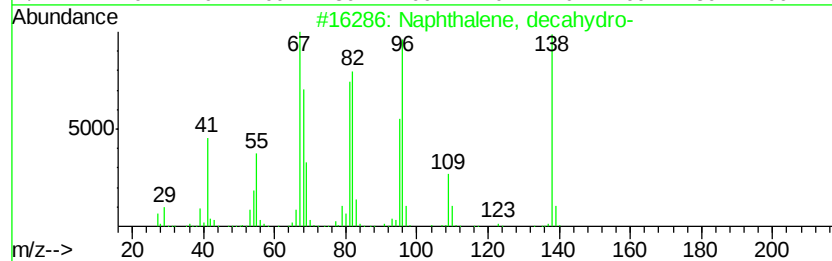
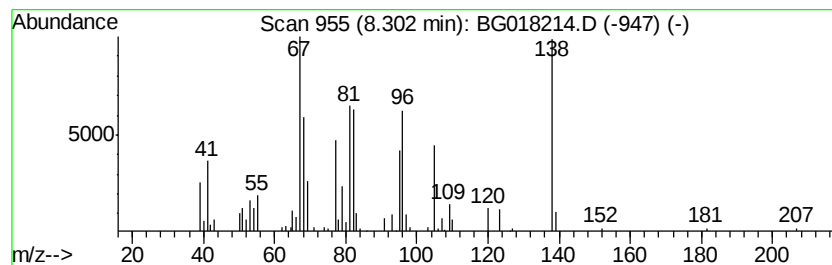
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Naphthalene, decahydro- Concentration Rank 29

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.30 | 5.65 ng/ul | 112584 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, decahydro-         | 138 | C10H18  | 000091-17-8 | 95   |
| 2    |    |   | Naphthalene, decahydro-         | 138 | C10H18  | 000091-17-8 | 95   |
| 3    |    |   | Naphthalene, decahydro-, trans- | 138 | C10H18  | 000493-02-7 | 93   |
| 4    |    |   | Naphthalene, decahydro-         | 138 | C10H18  | 000091-17-8 | 91   |
| 5    |    |   | Naphthalene, decahydro-, cis-   | 138 | C10H18  | 000493-01-6 | 90   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

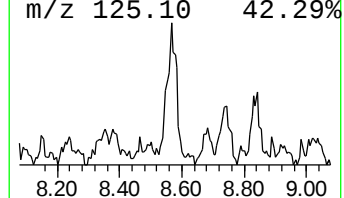
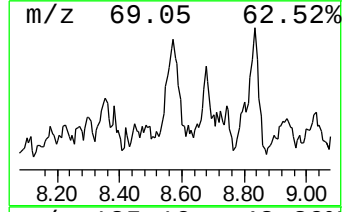
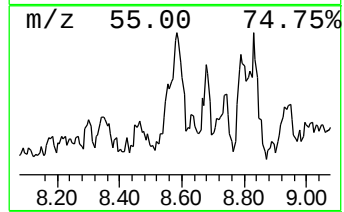
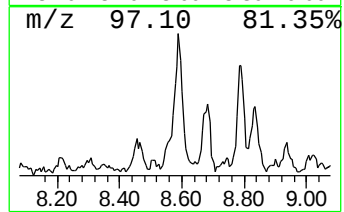
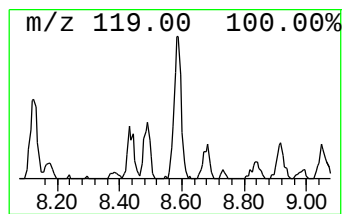
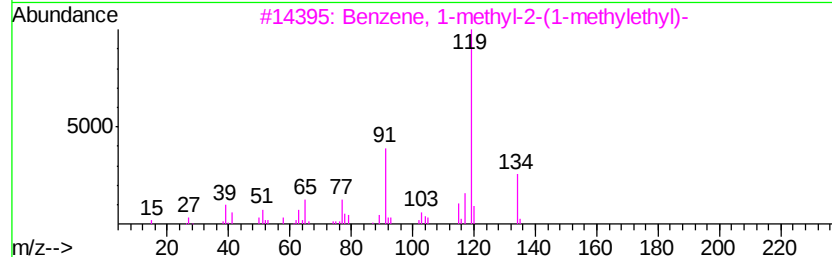
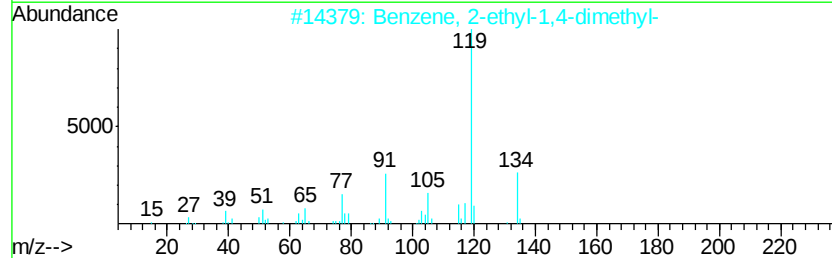
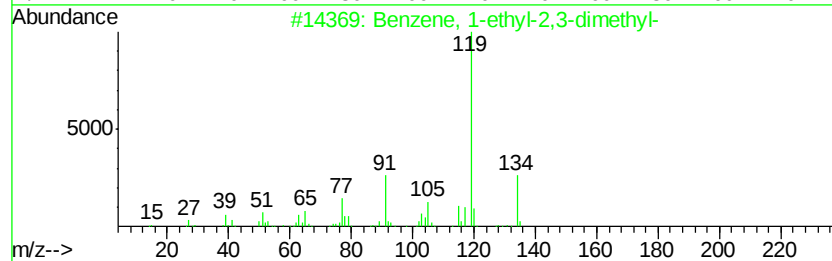
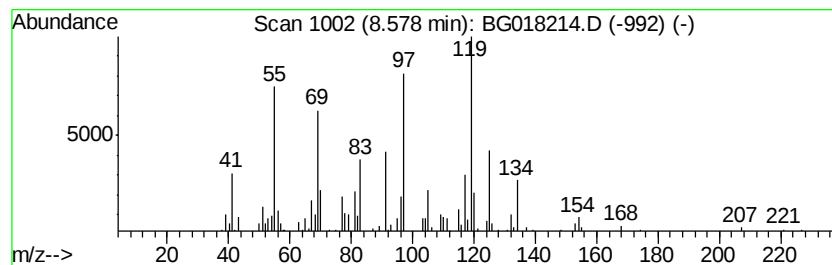
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 17

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.58 | 9.67 ng/ul | 192711 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 86   |
| 2    |    | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 46   |
| 3    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 46   |
| 4    |    | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 46   |
| 5    |    | Benzene, 1-methyl-4-(1-methyleth... | 134 | C10H14  | 000099-87-6 | 46   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

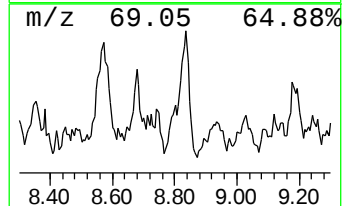
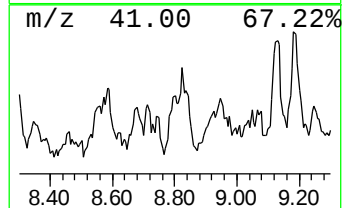
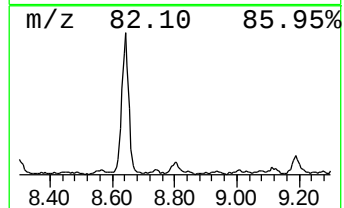
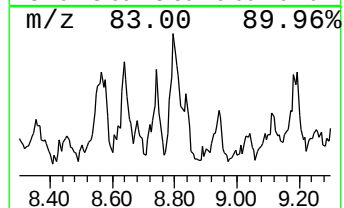
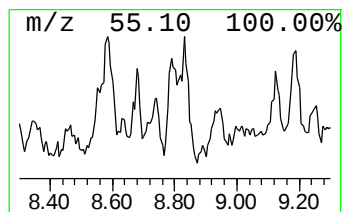
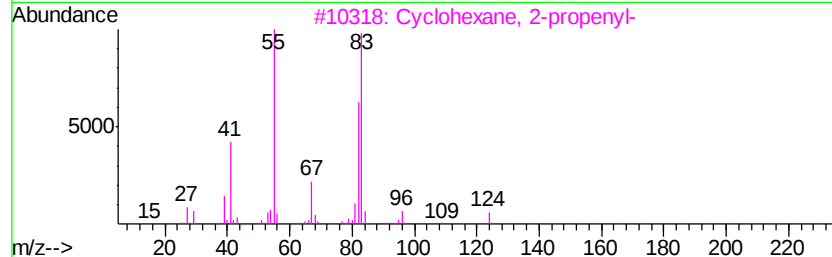
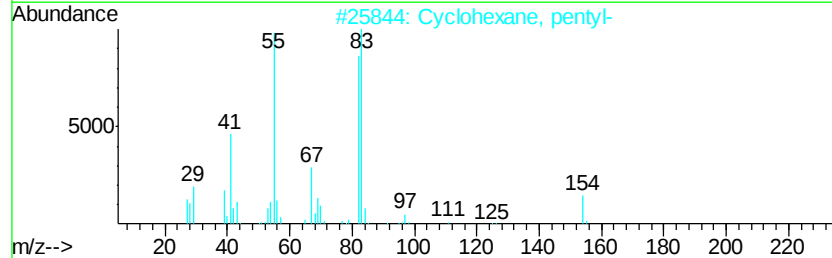
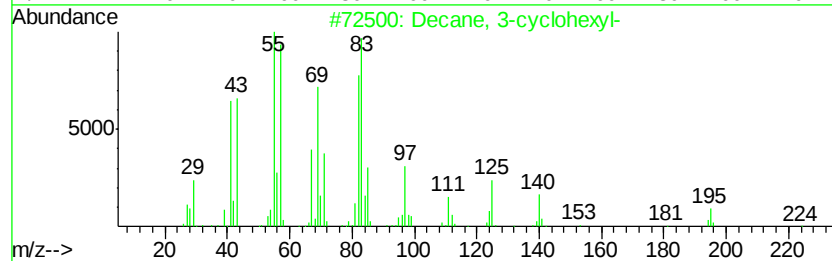
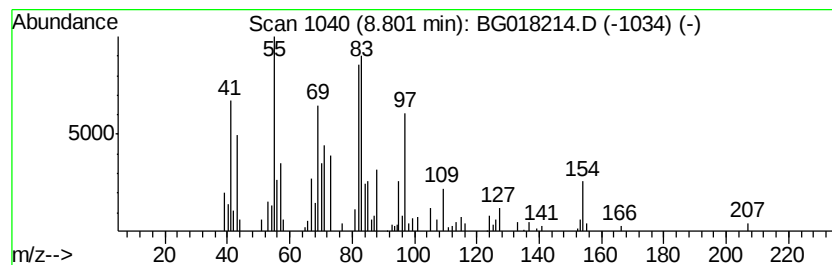
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 38

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 8.80 | 4.32 ng/ul | 86032 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 3-cyclohexyl-           | 224 | C16H32  | 013151-74-1 | 53   |
| 2    |    | Cyclohexane, pentyl-            | 154 | C11H22  | 004292-92-6 | 43   |
| 3    |    | Cyclohexane, 2-propenyl-        | 124 | C9H16   | 002114-42-3 | 38   |
| 4    |    | Cyclohexane, 1,1'-methylenebis- | 180 | C13H24  | 003178-23-2 | 38   |
| 5    |    | Cyclohexane, (1-methylethyl)-   | 126 | C9H18   | 000696-29-7 | 38   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

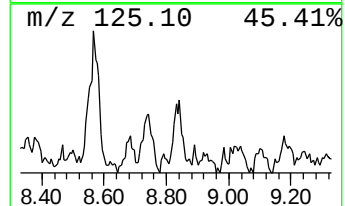
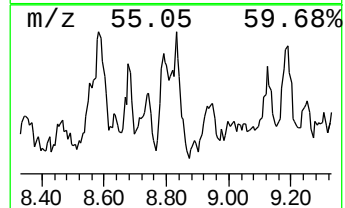
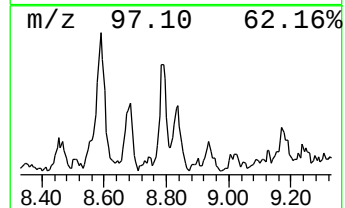
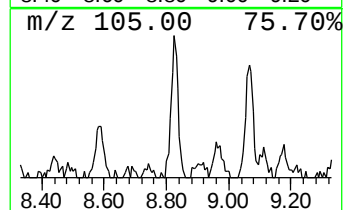
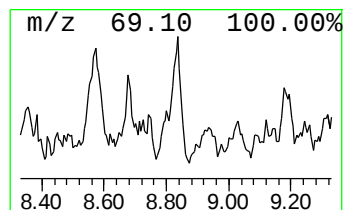
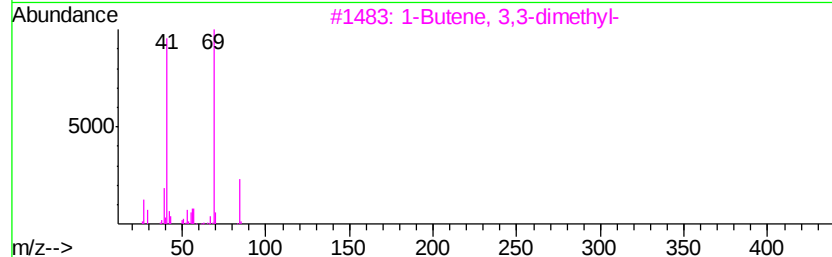
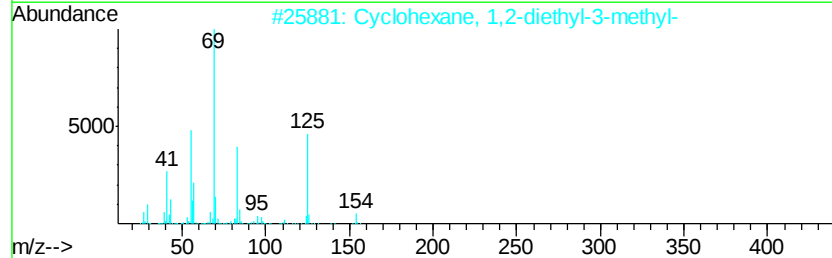
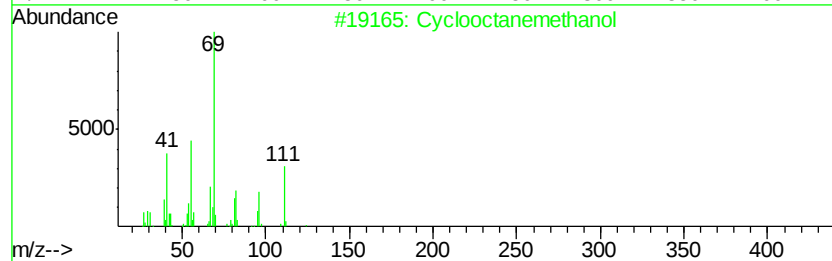
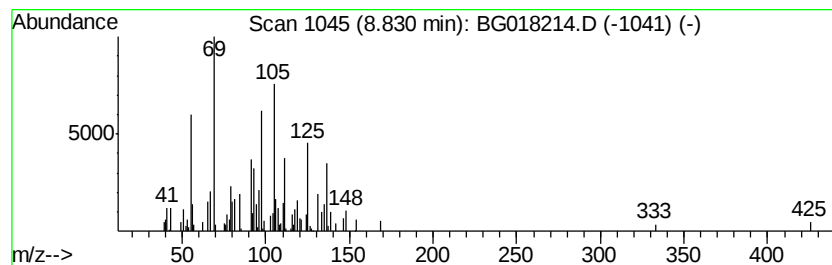
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 unknown-03 Concentration Rank 23

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.83 | 7.95 ng/ul | 158518 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclooctanemethanol                 | 142 | C9H18O  | 003637-63-6 | 35   |
| 2    |    | Cyclohexane, 1,2-diethyl-3-methyl-  | 154 | C11H22  | 061141-80-8 | 35   |
| 3    |    | 1-Butene, 3,3-dimethyl-             | 84  | C6H12   | 000558-37-2 | 25   |
| 4    |    | Cyclopentane, 1,1,3,4-tetramethy... | 126 | C9H18   | 053907-60-1 | 25   |
| 5    |    | Cyclohexane, 1-ethyl-2-propyl-      | 154 | C11H22  | 062238-33-9 | 25   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

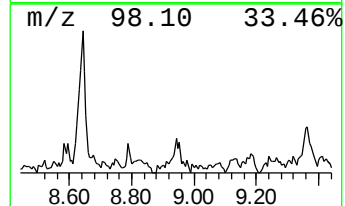
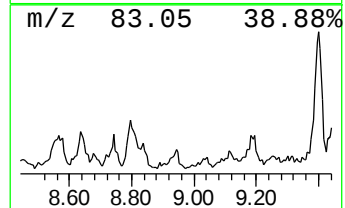
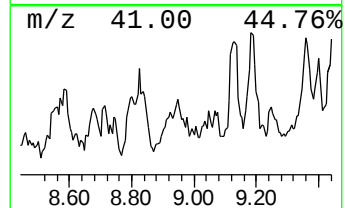
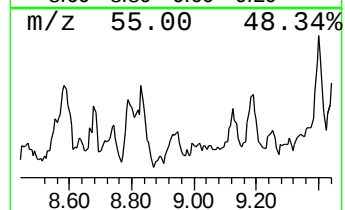
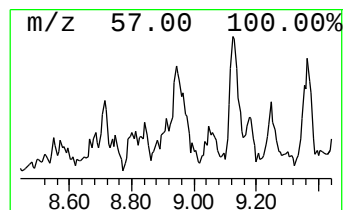
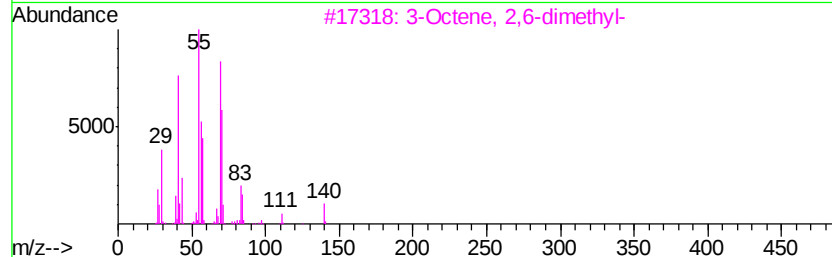
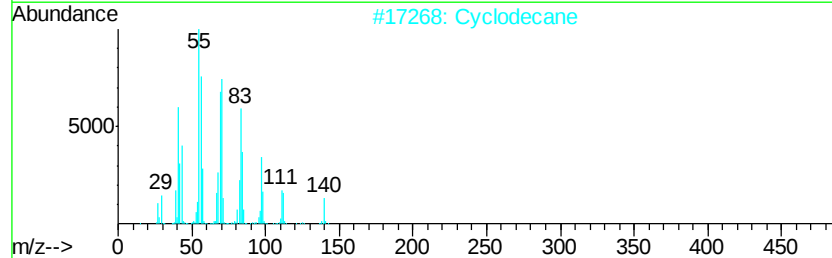
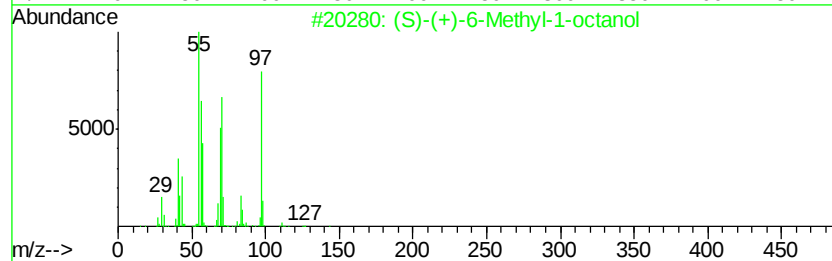
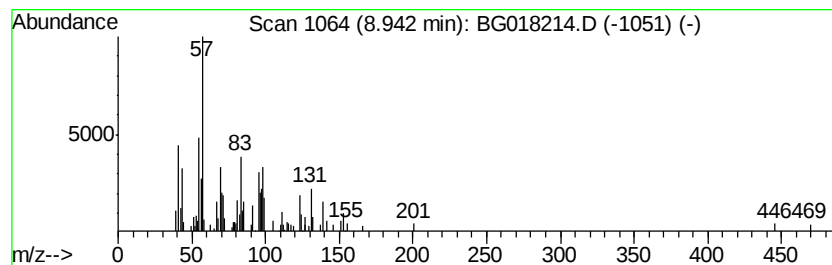
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 unknown-04 Concentration Rank 37

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 8.94 | 4.44 ng/ul | 170899 | Naphthalene-d8   | 10.27 |

| Hit# | of                                     | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----------------------------------------|---|--------------|-----|---------|-------------|------|
| 1    | (S)-(+)-6-Methyl-1-octanol             |   |              | 144 | C9H20O  | 110453-78-6 | 27   |
| 2    | Cyclodecane                            |   |              | 140 | C10H20  | 000293-96-9 | 18   |
| 3    | 3-Octene, 2,6-dimethyl-                |   |              | 140 | C10H20  | 006874-28-8 | 15   |
| 4    | Cyclohexane, 1-methyl-4-(1-methyl-...) |   |              | 140 | C10H20  | 001678-82-6 | 14   |
| 5    | Cyclopropane, 1-ethyl-2-heptyl-        |   |              | 168 | C12H24  | 074663-86-8 | 14   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

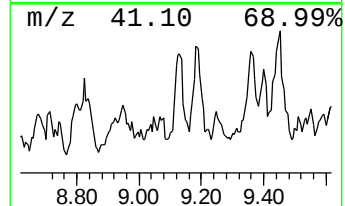
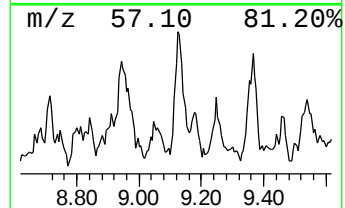
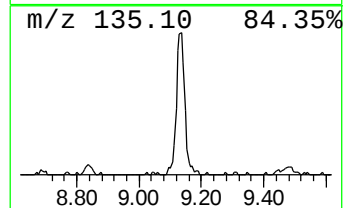
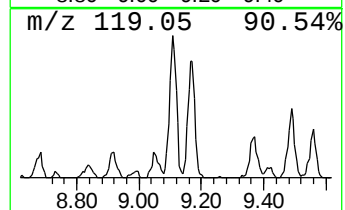
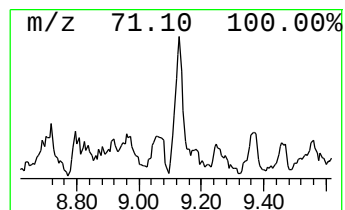
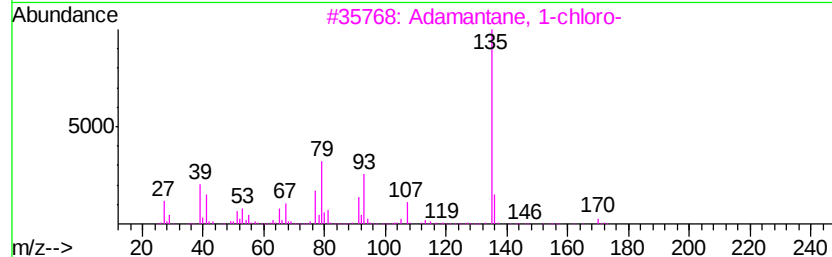
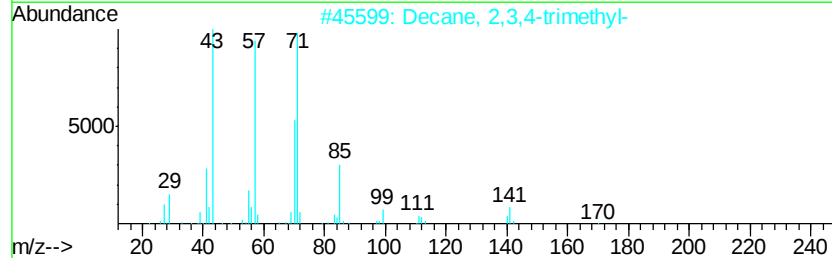
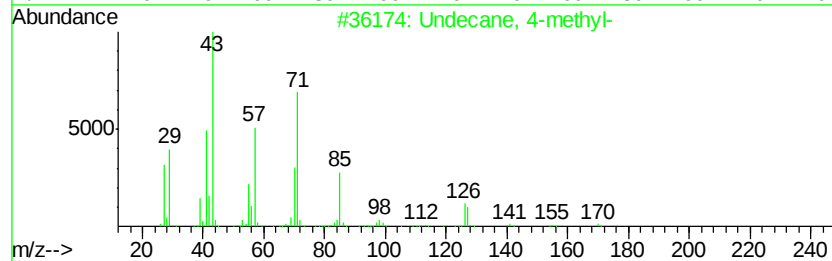
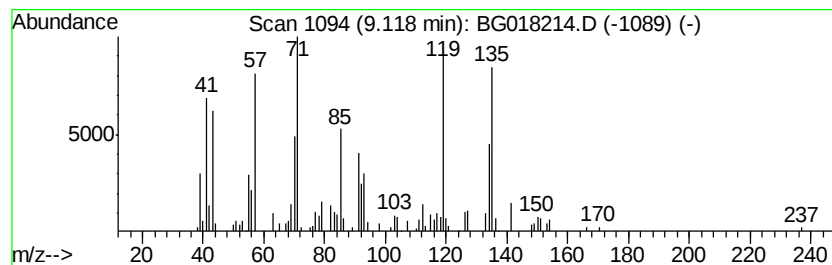
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.12 | 4.04 ng/ul | 155388 | Naphthalene-d8   | 10.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Undecane, 4-methyl-                 | 170 | C12H26    | 002980-69-0  | 38   |
| 2    |    | Decane, 2,3,4-trimethyl-            | 184 | C13H28    | 062238-15-7  | 27   |
| 3    |    | Adamantane, 1-chloro-               | 170 | C10H15Cl  | 000935-56-8  | 25   |
| 4    |    | m-Aminobenzamidine                  | 135 | C7H9N3    | 003459-66-3  | 22   |
| 5    |    | 4-Cyanobenzoic acid, 2-(1-adaman... | 309 | C20H23NO2 | 1000292-45-1 | 22   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

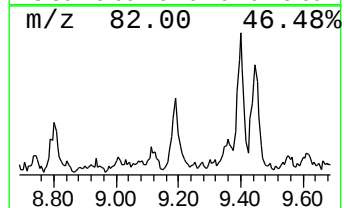
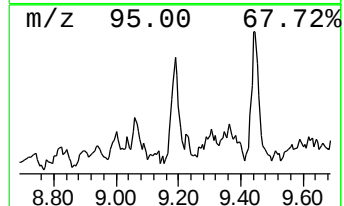
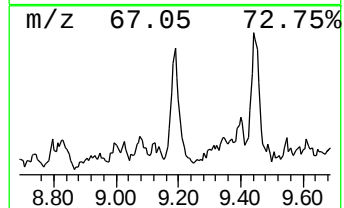
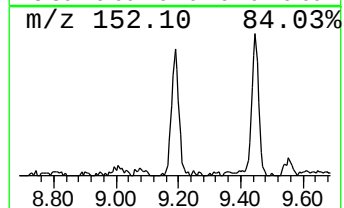
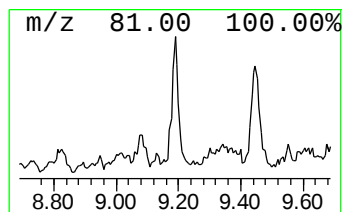
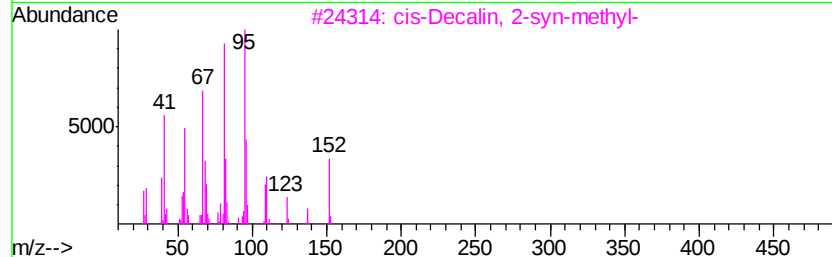
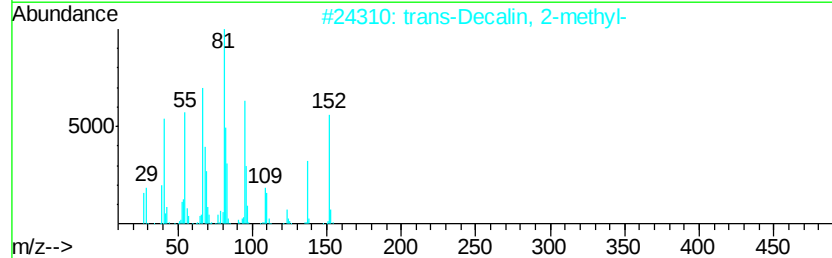
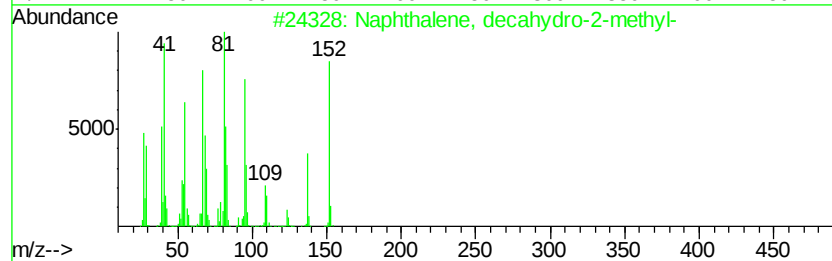
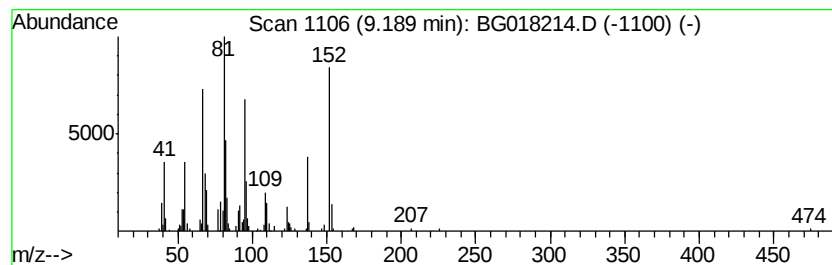
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Naphthalene, decahydro-2-me... Concentration Rank 31

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.19 | 5.11 ng/ul | 196404 | Naphthalene-d8   | 10.27 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Naphthalene, decahydro-2-methyl-    | 152 | C11H20  | 002958-76-1  | 91   |
| 2    |    |   | trans-Decalin, 2-methyl-            | 152 | C11H20  | 1000152-47-3 | 91   |
| 3    |    |   | cis-Decalin, 2-syn-methyl-          | 152 | C11H20  | 1000155-85-6 | 91   |
| 4    |    |   | Cyclohexanone, 5-methyl-2-(1-met... | 152 | C10H16O | 015932-80-6  | 90   |
| 5    |    |   | Pulegone                            | 152 | C10H16O | 000089-82-7  | 81   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

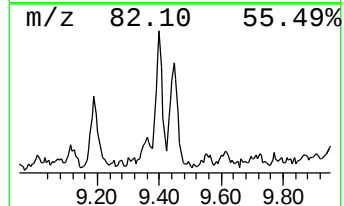
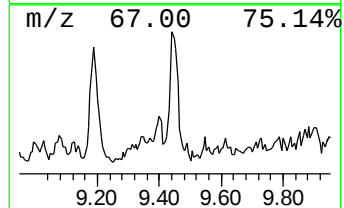
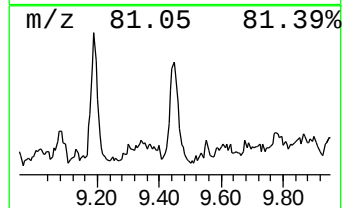
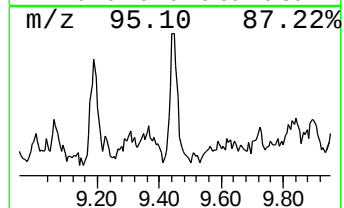
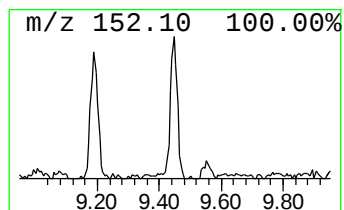
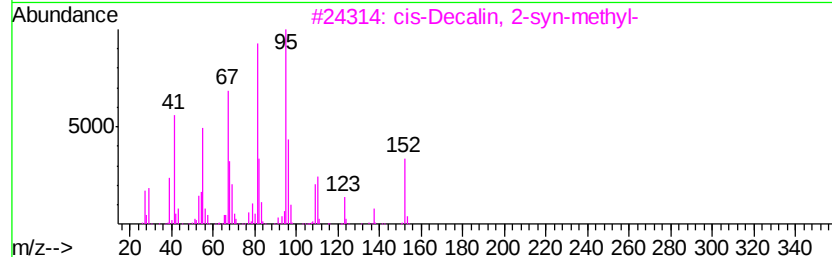
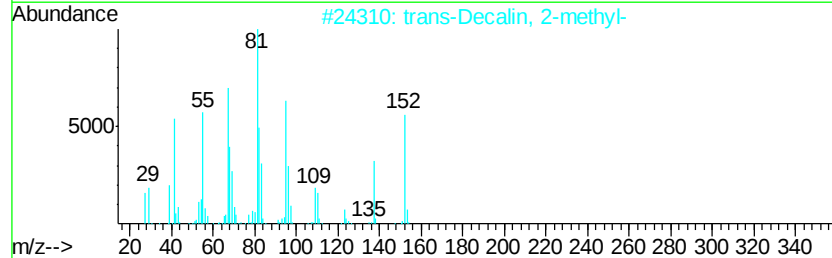
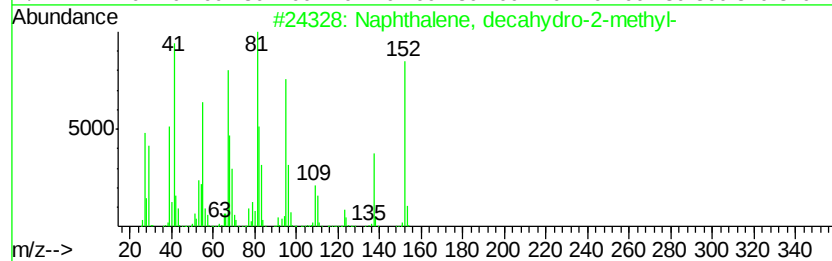
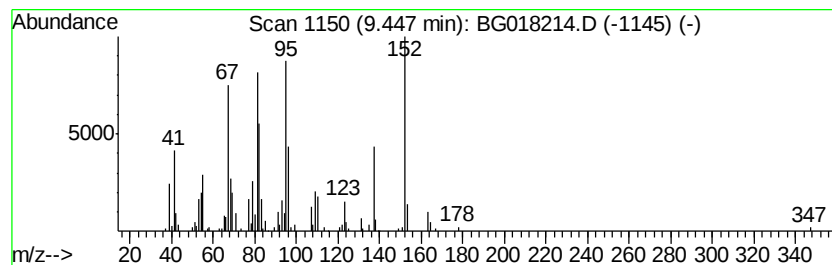
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 trans-Decalin, 2-methyl- Concentration Rank 27

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.45 | 6.00 ng/ul | 230631 | Naphthalene-d8   | 10.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Naphthalene, decahydro-2-methyl-    | 152 | C11H20  | 002958-76-1  | 90   |
| 2    |    | trans-Decalin, 2-methyl-            | 152 | C11H20  | 1000152-47-3 | 87   |
| 3    |    | cis-Decalin, 2-syn-methyl-          | 152 | C11H20  | 1000155-85-6 | 80   |
| 4    |    | Naphthalene, decahydro-2-methyl-    | 152 | C11H20  | 002958-76-1  | 64   |
| 5    |    | 1,3,3-Trimethylcyclohex-1-ene-4-... | 152 | C10H16O | 127128-60-3  | 52   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

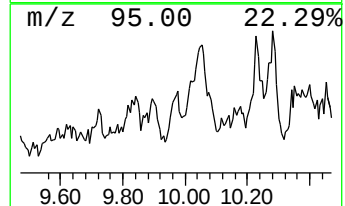
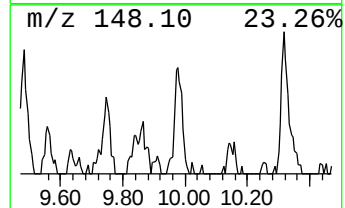
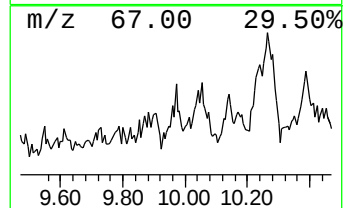
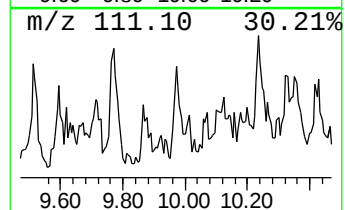
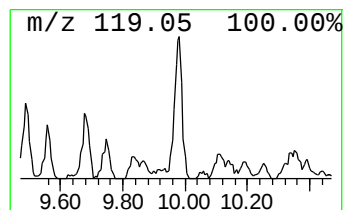
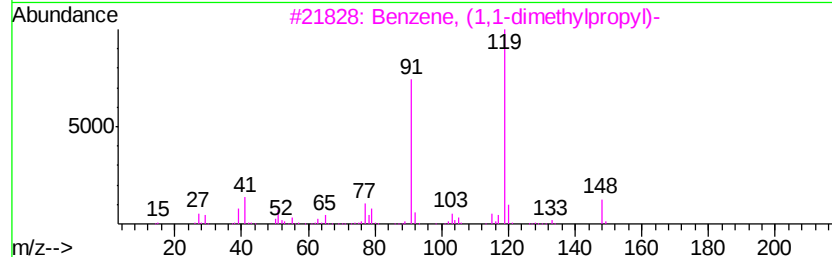
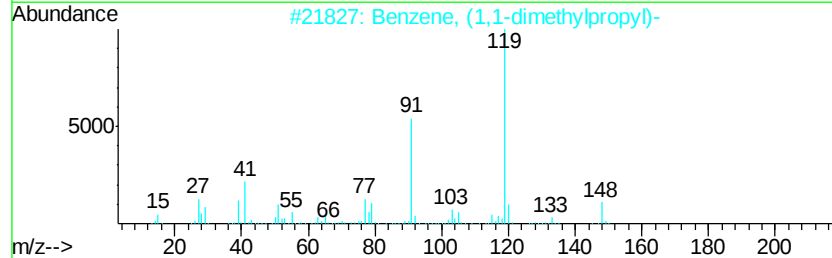
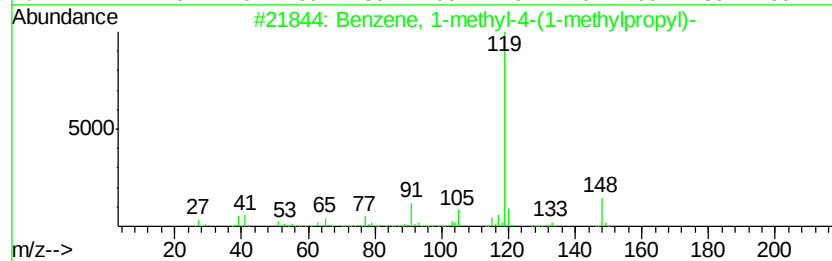
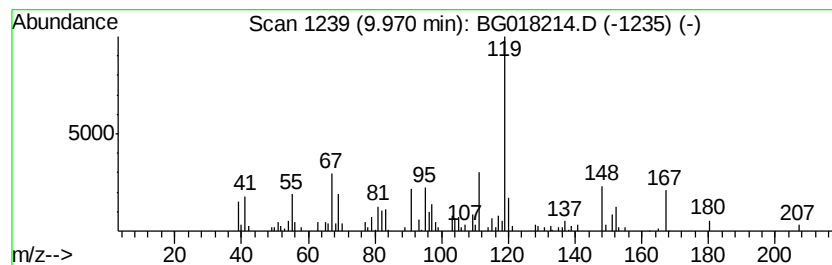
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Benzene, 1-methyl-4-(1-meth... Concentration Rank 56

| R.T. | EstConc    | Area  | Relative to ISTD | R.T.  |
|------|------------|-------|------------------|-------|
| 9.97 | 2.23 ng/ul | 85661 | Napthalene-d8    | 10.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16   | 001595-16-0  | 50   |
| 2    |    | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16   | 002049-95-8  | 38   |
| 3    |    | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16   | 002049-95-8  | 38   |
| 4    |    | o-Toluylic acid, oct-3-en-2-yl e... | 246 | C16H22O2 | 1000292-51-3 | 35   |
| 5    |    | Benzene, 2-ethyl-1,3-dimethyl-      | 134 | C10H14   | 002870-04-4  | 27   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

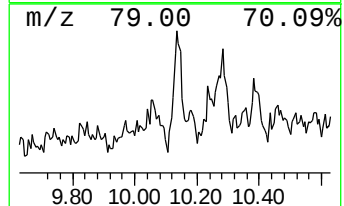
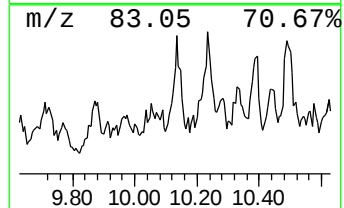
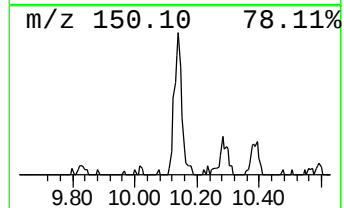
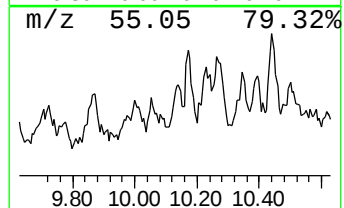
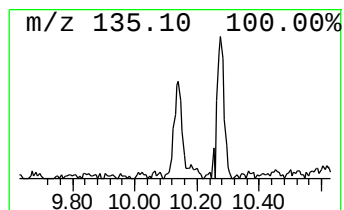
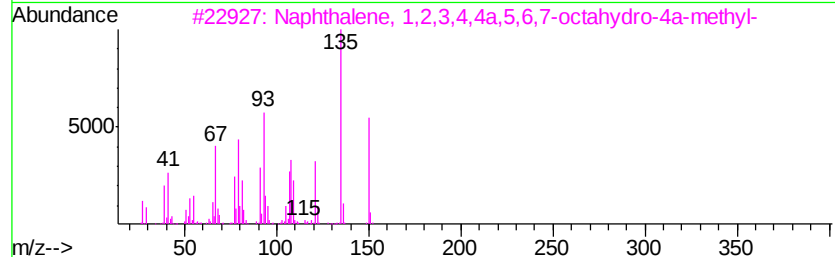
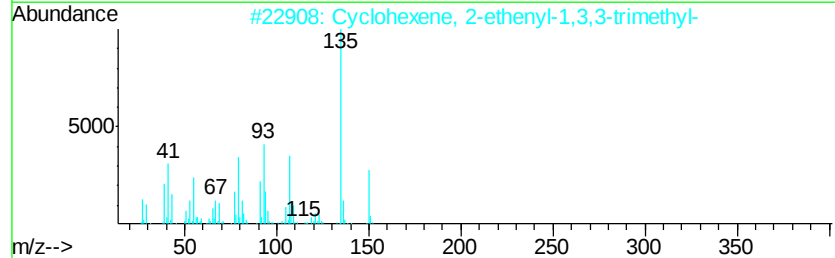
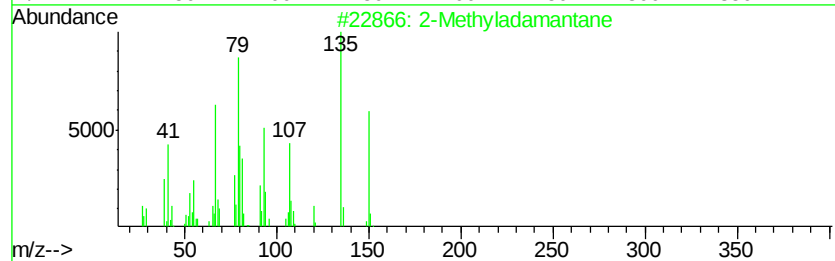
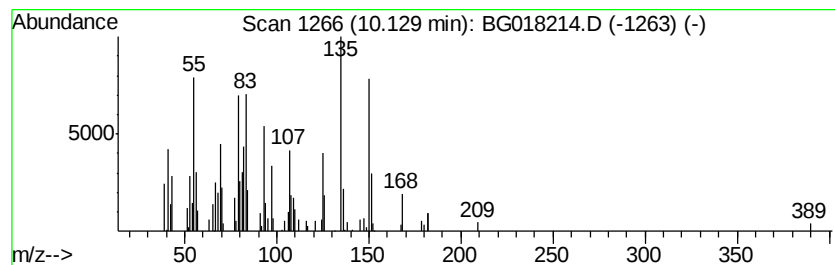
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 2-Methyladamantane Concentration Rank 55

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 10.13 | 2.46 ng/ul | 94770 | Naphthalene-d8   | 10.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Methyladamantane                  | 150 | C11H18  | 000700-56-1 | 60   |
| 2    |    | Cyclohexene, 2-ethenyl-1,3,3-tri... | 150 | C11H18  | 005293-90-3 | 49   |
| 3    |    | Naphthalene, 1,2,3,4,4a,5,6,7-oc... | 150 | C11H18  | 013943-77-6 | 49   |
| 4    |    | 2H-Inden-2-one, 1,4,5,6,7,7a-hex... | 150 | C10H14O | 054725-16-5 | 46   |
| 5    |    | Cyclohexanone, 2-(2-butyryl)-       | 150 | C10H14O | 054166-48-2 | 46   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

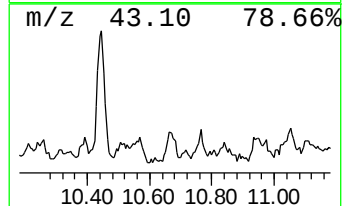
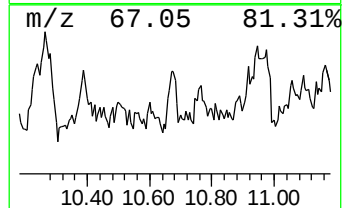
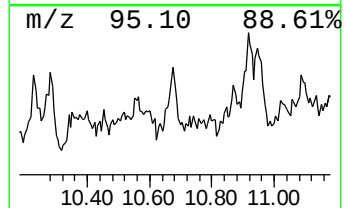
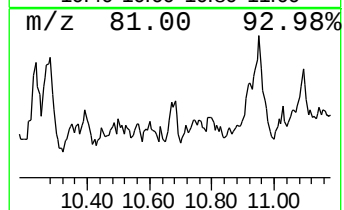
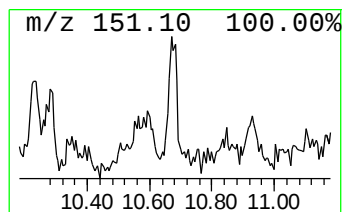
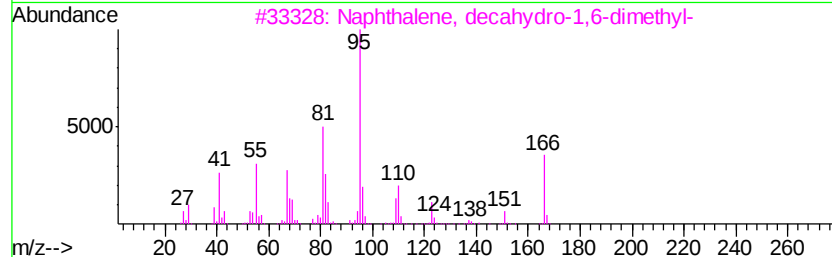
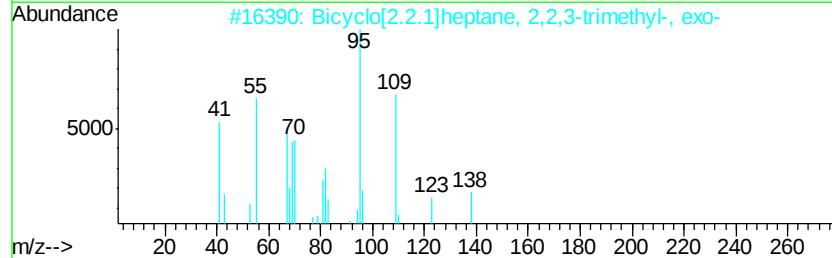
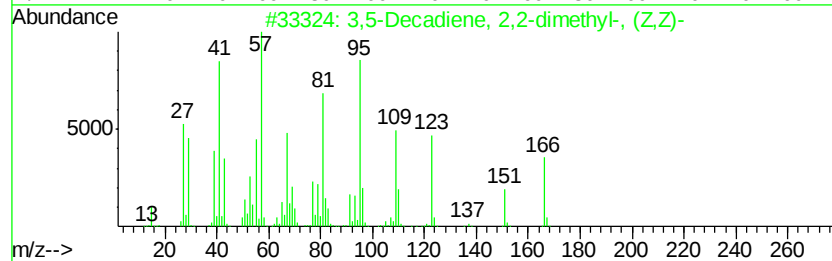
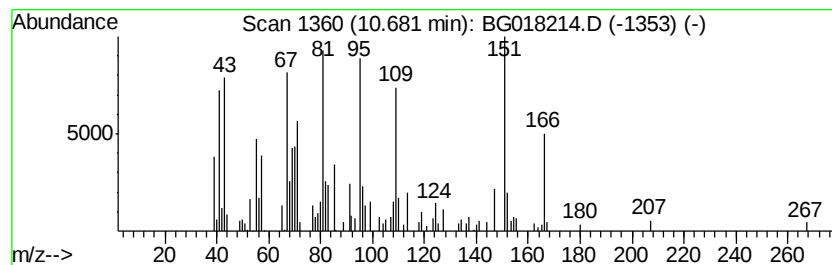
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 unknown-05 Concentration Rank 49

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.68 | 2.99 ng/ul | 114975 | Naphthalene-d8   | 10.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 3,5-Decadiene, 2,2-dimethyl-, (Z... | 166 | C12H22  | 055638-50-1 | 43   |
| 2    |    | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138 | C10H18  | 020536-41-8 | 42   |
| 3    |    | Naphthalene, decahydro-1,6-dimet... | 166 | C12H22  | 001750-51-2 | 42   |
| 4    |    | Tricyclo[4.3.1.1(3,8)]undecan-3-ol  | 166 | C11H18O | 014504-80-4 | 41   |
| 5    |    | Naphthalene, decahydro-2,3-dimet... | 166 | C12H22  | 001008-80-6 | 41   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

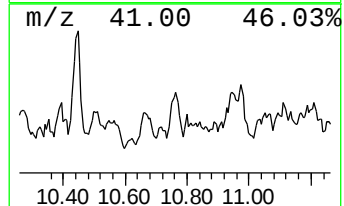
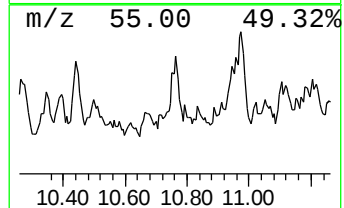
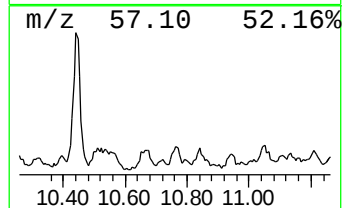
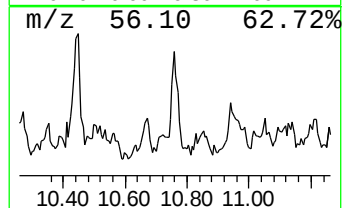
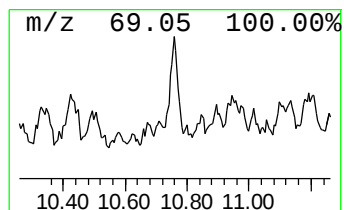
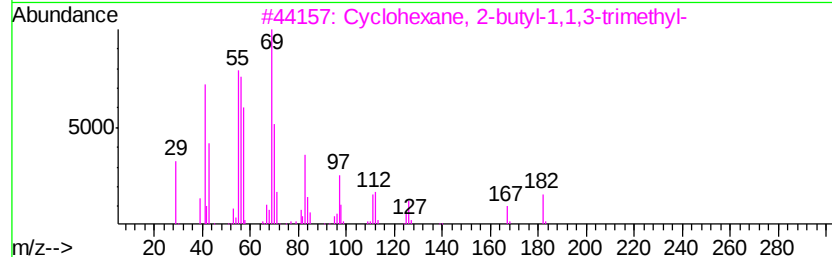
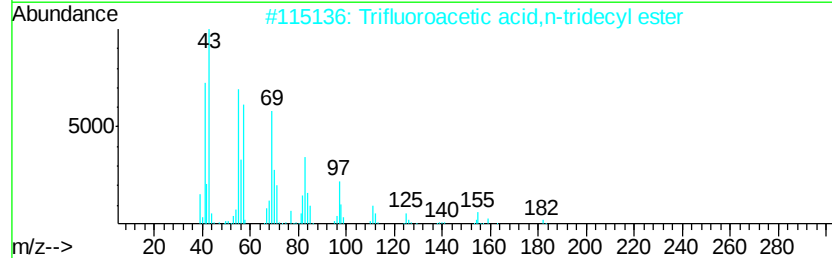
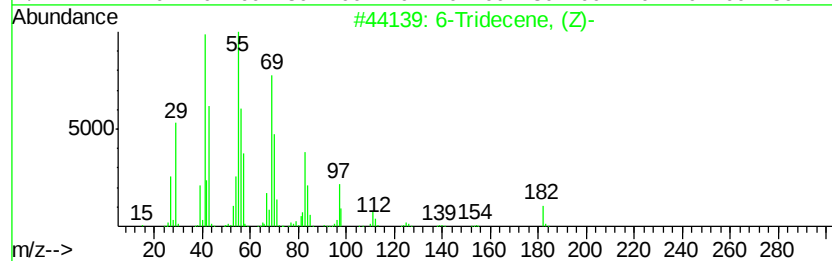
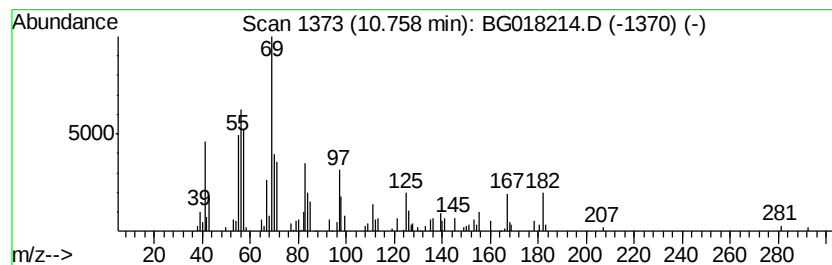
Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 18 6-Tridecene, (Z)- Concentration Rank 53

| R.T.    | EstConc                             | Area           | Relative to ISTD | R.T.      |
|---------|-------------------------------------|----------------|------------------|-----------|
| 10.76   | 2.55 ng/ul                          | 98004          | Naphthalene-d8   | 10.27     |
| Hit# of | 5                                   | Tentative ID   | MW MolForm       | CAS# Qual |
| 1       | 6-Tridecene, (Z)-                   | 182 C13H26     | 006508-77-6      | 64        |
| 2       | Trifluoroacetic acid,n-tridecyl ... | 296 C15H27F3O2 | 053800-02-5      | 62        |
| 3       | Cyclohexane, 2-butyl-1,1,3-trime... | 182 C13H26     | 054676-39-0      | 60        |
| 4       | 4-Tridecene, (Z)-                   | 182 C13H26     | 041446-54-2      | 55        |
| 5       | Cyclohexane, 2-butyl-1,1,3-trime... | 182 C13H26     | 054676-39-0      | 53        |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

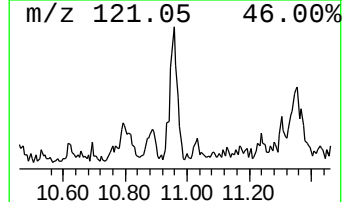
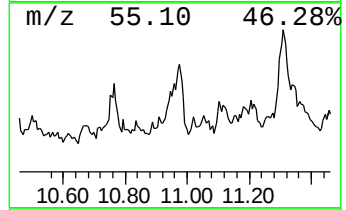
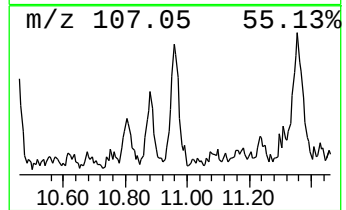
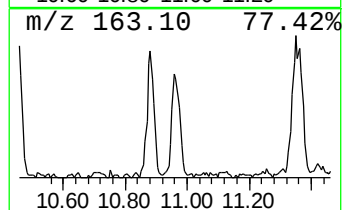
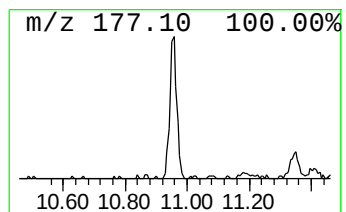
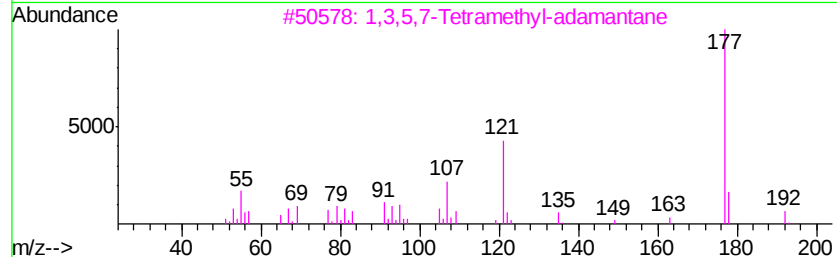
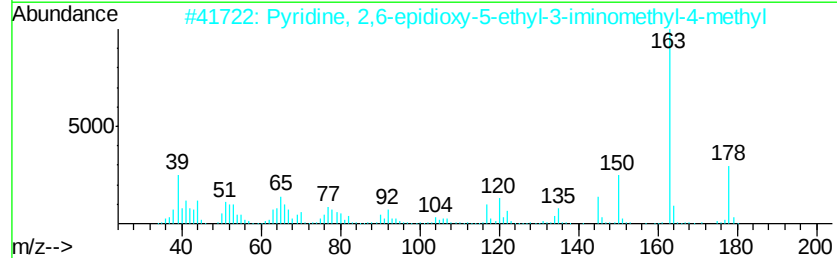
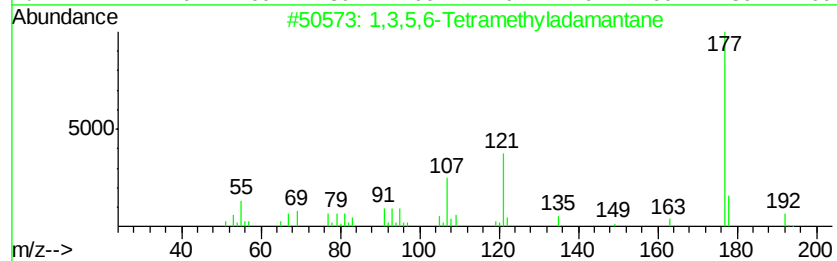
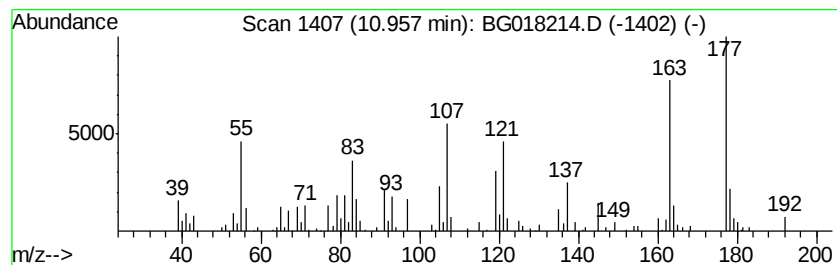
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 unknown-06 Concentration Rank 20

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.96 | 8.79 ng/ul | 338196 | Naphthalene-d8   | 10.27 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 1,3,5,6-Tetramethyladamantane       | 192 | C14H24    | 034694-70-7  | 49   |
| 2    |    | Pyridine, 2,6-epidioxy-5-ethyl-3... | 178 | C9H10N2O2 | 1000263-31-1 | 41   |
| 3    |    | 1,3,5,7-Tetramethyl-adamantane      | 192 | C14H24    | 001687-36-1  | 41   |
| 4    |    | 1,3-Disilaindan, 1,3-dimethyl-      | 178 | C9H14Si2  | 017864-73-2  | 35   |
| 5    |    | Phenol, 2,4-bis(1-methylethyl)-     | 178 | C12H18O   | 002934-05-6  | 35   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

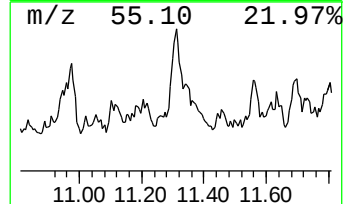
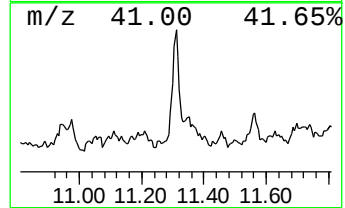
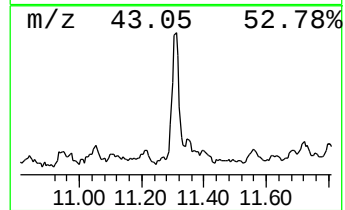
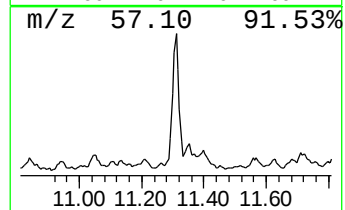
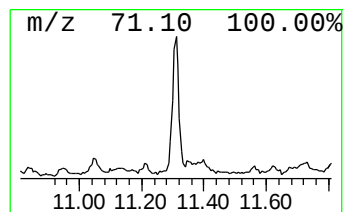
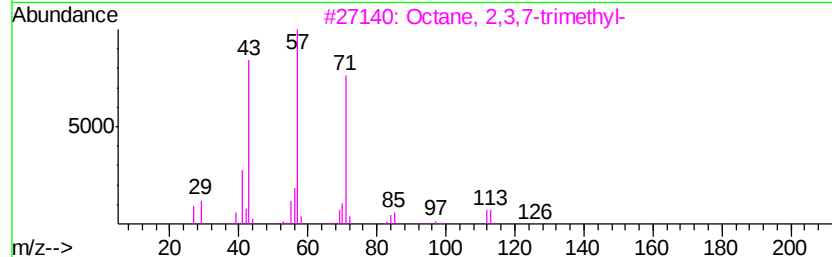
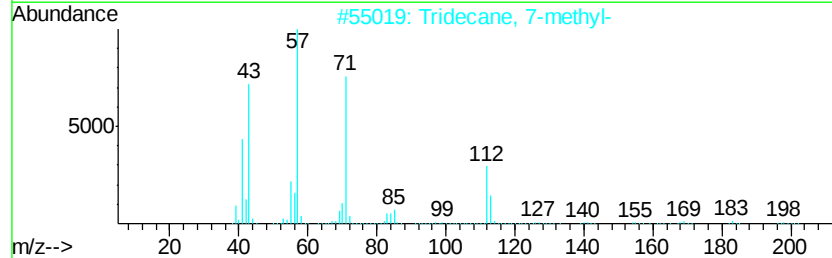
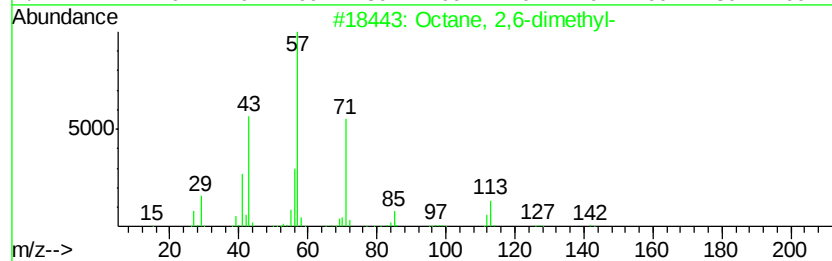
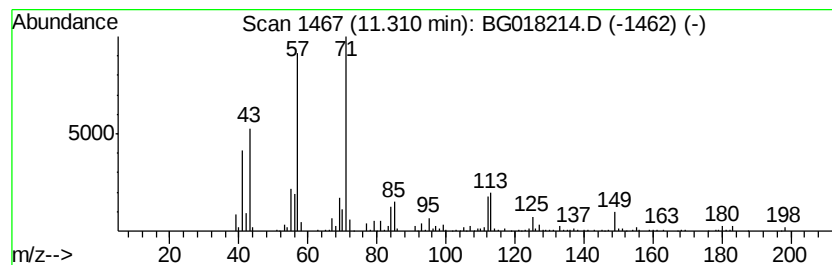
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 11.31 | 13.94 ng/ul | 536213 | Naphthalene-d8   | 10.27 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,6-dimethyl-    | 142 | C10H22  | 002051-30-1 | 80   |
| 2    |    |   | Tridecane, 7-methyl-     | 198 | C14H30  | 026730-14-3 | 76   |
| 3    |    |   | Octane, 2,3,7-trimethyl- | 156 | C11H24  | 062016-34-6 | 72   |
| 4    |    |   | Decane, 4-methyl-        | 156 | C11H24  | 002847-72-5 | 59   |
| 5    |    |   | Octane, 2-bromo-         | 192 | C8H17Br | 000557-35-7 | 50   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

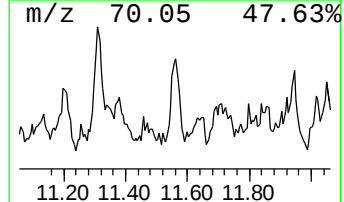
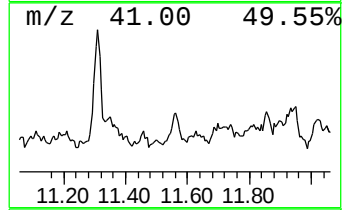
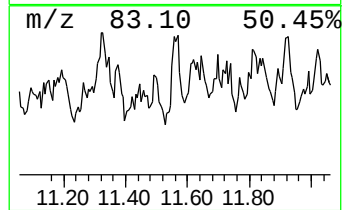
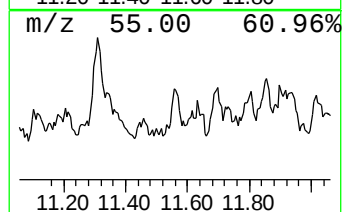
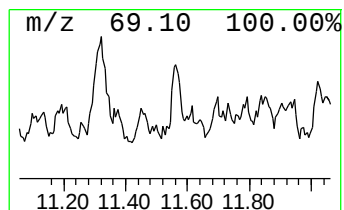
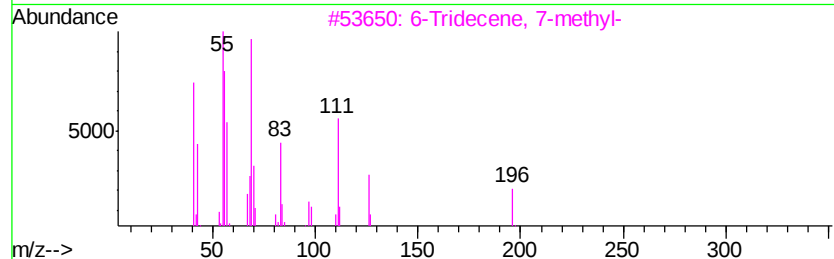
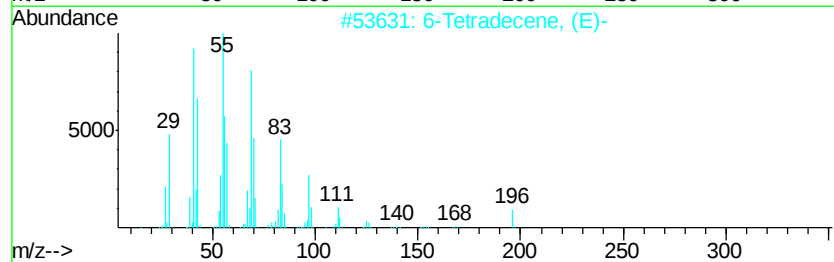
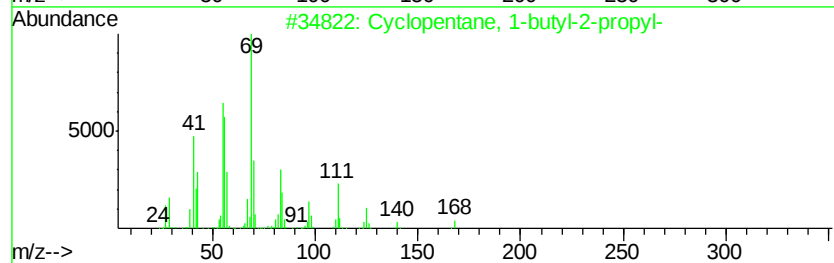
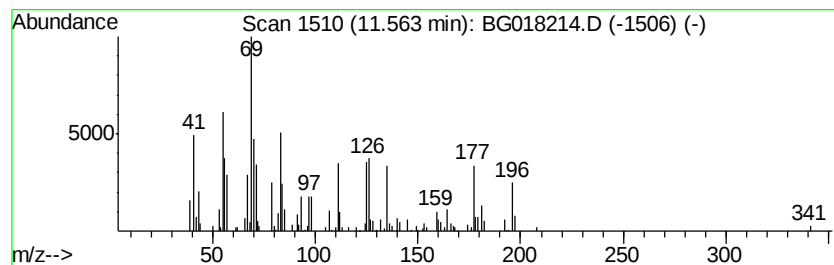
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 47

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.56 | 3.15 ng/ul | 120979 | Naphthalene-d8   | 10.27 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclopentane, 1-butyl-2-propyl- | 168 | C12H24  | 062199-50-2 | 70   |
| 2    |    |   | 6-Tetradecene, (E)-             | 196 | C14H28  | 041446-64-4 | 62   |
| 3    |    |   | 6-Tridecene, 7-methyl-          | 196 | C14H28  | 024949-42-6 | 62   |
| 4    |    |   | Cyclopentane, 1-butyl-2-pentyl- | 196 | C14H28  | 061142-52-7 | 60   |
| 5    |    |   | Cyclotetradecane                | 196 | C14H28  | 000295-17-0 | 46   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

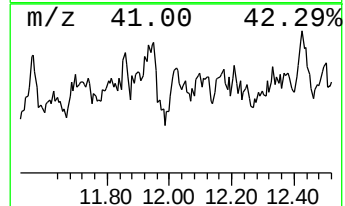
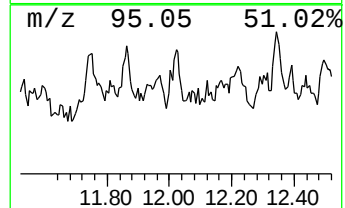
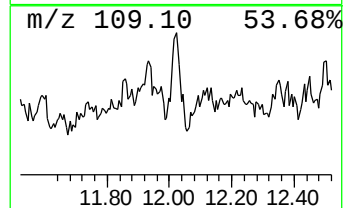
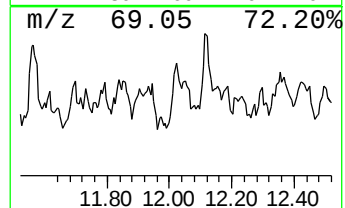
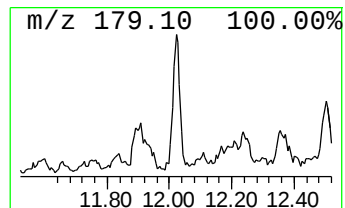
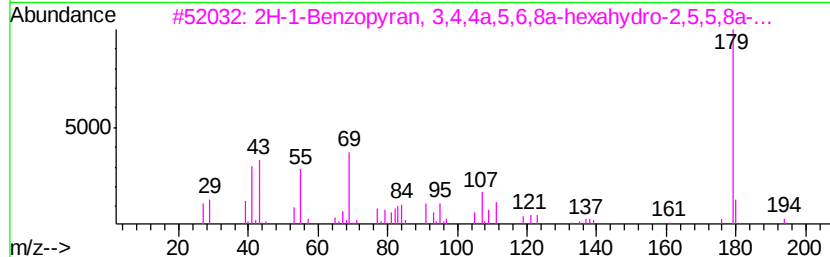
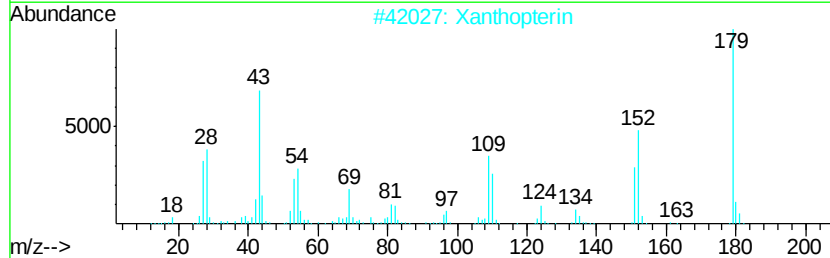
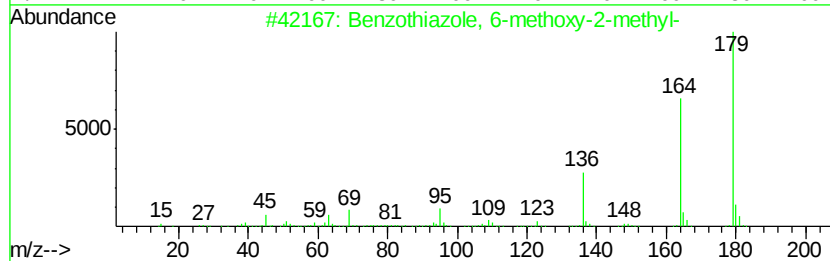
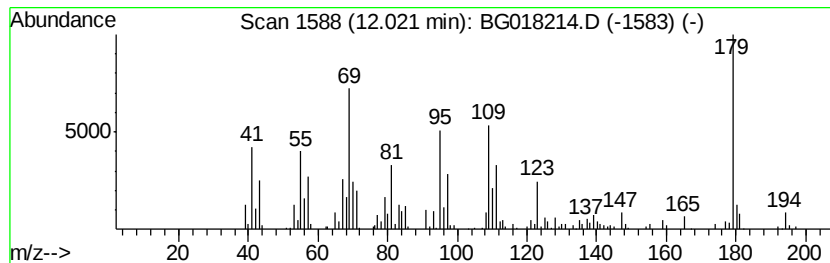
Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 22 unknown-07 Concentration Rank 34

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.           |
|---------|------------|-------------------------------------|------------------|----------------|
| 12.02   | 4.75 ng/ul | 182775                              | Napthalene-d8    | 10.27          |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |            | Benzothiazole, 6-methoxy-2-methyl-  | 179 C9H9NOS      | 002941-72-2 43 |
| 2       |            | Xanthopterin                        | 179 C6H5N5O2     | 005979-01-1 38 |
| 3       |            | 2H-1-Benzopyran, 3,4,4a,5,6,8a-h... | 194 C13H22O      | 063335-66-0 35 |
| 4       |            | 2(3H)-Benzothiazolone, 3-methyl-... | 179 C8H9N3S      | 001128-67-2 27 |
| 5       |            | 2,4-Dimethoxyphenyl isocyanate      | 179 C9H9NO3      | 084370-87-6 25 |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

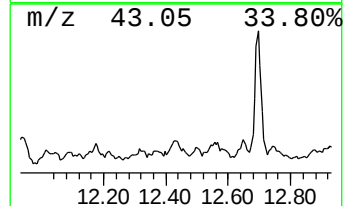
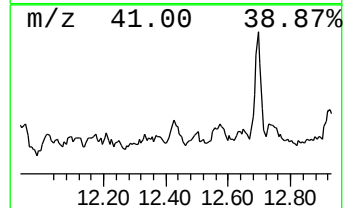
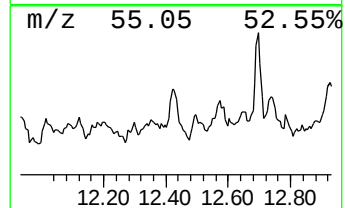
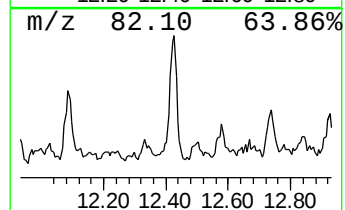
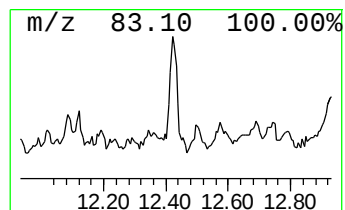
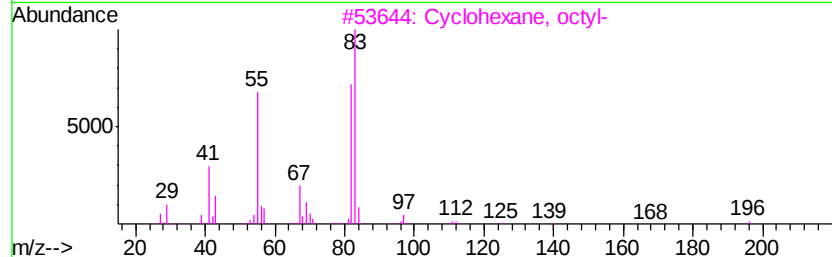
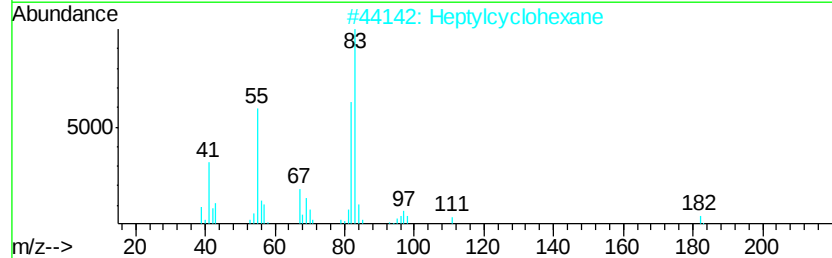
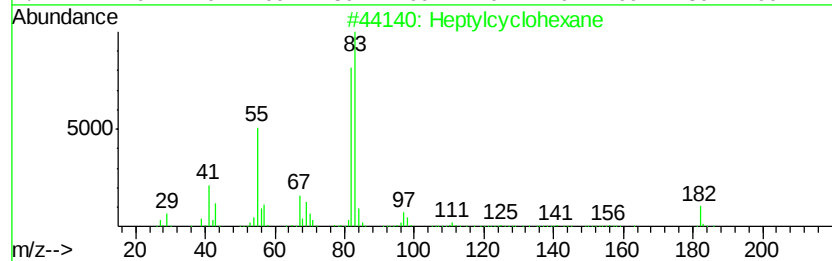
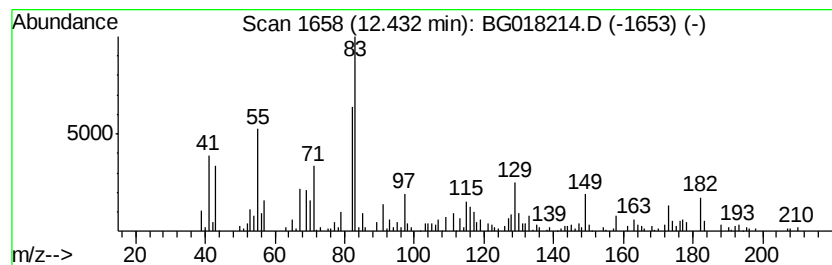
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 (DEL) Alkane: Cyclic12.43 Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.43 | 5.71 ng/ul | 199506 | Acenaphthene-d10 | 14.17 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Heptylcyclohexane    | 182 | C13H26  | 005617-41-4 | 49   |
| 2    |    |   | Heptylcyclohexane    | 182 | C13H26  | 005617-41-4 | 47   |
| 3    |    |   | Cyclohexane, octyl-  | 196 | C14H28  | 001795-15-9 | 46   |
| 4    |    |   | Cyclohexane, pentyl- | 154 | C11H22  | 004292-92-6 | 43   |
| 5    |    |   | Cyclohexane, octyl-  | 196 | C14H28  | 001795-15-9 | 43   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

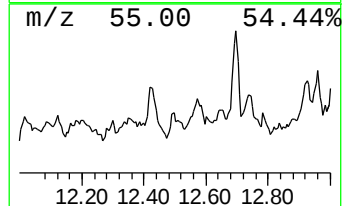
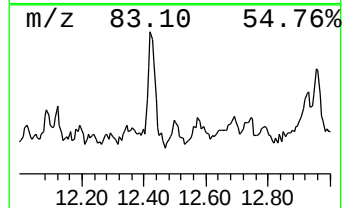
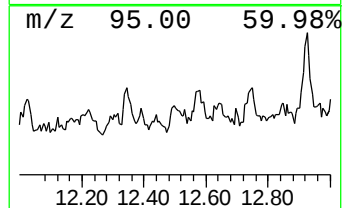
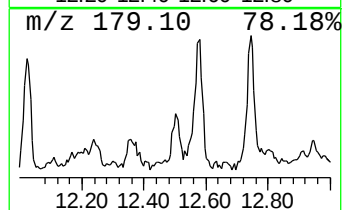
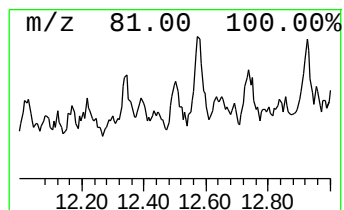
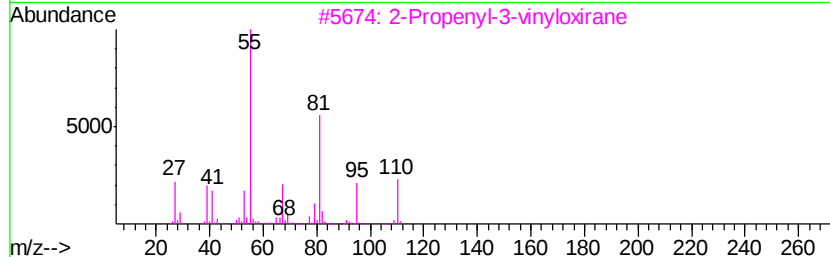
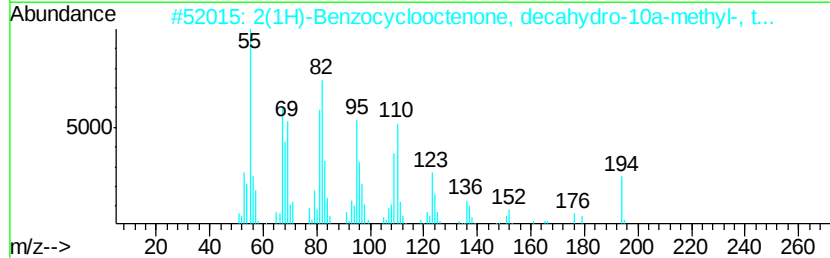
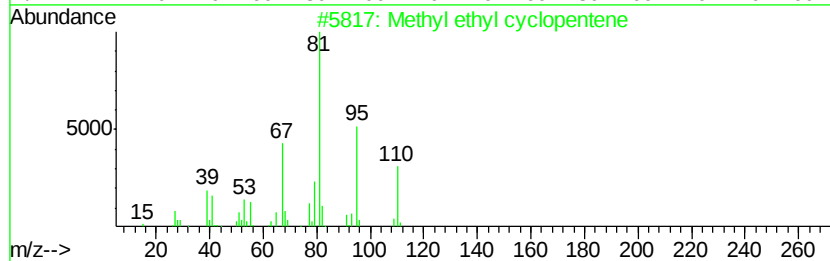
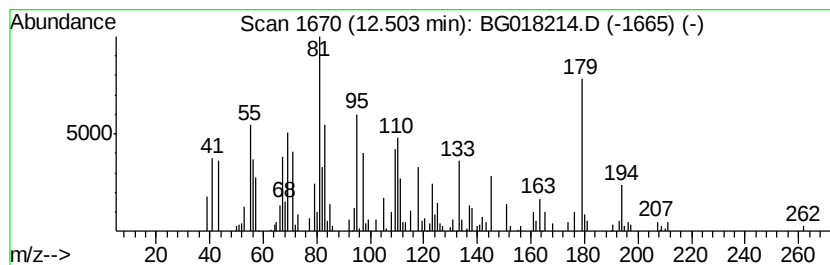
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 unknown-08 Concentration Rank 42

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.50 | 3.81 ng/ul | 133177 | Acenaphthene-d10 | 14.17 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Methyl ethyl cyclopentene           | 110 | C8H14   | 019780-56-4 | 30   |
| 2    |    |   | 2(1H)-Benzocyclooctenone, decahy... | 194 | C13H22O | 055103-68-9 | 30   |
| 3    |    |   | 2-Propenyl-3-vinylloxirane          | 110 | C7H10O  | 070597-49-8 | 27   |
| 4    |    |   | 1-Ethyl-5-methylcyclopentene        | 110 | C8H14   | 097797-57-4 | 27   |
| 5    |    |   | Cyclohexane, ethenyl-               | 110 | C8H14   | 000695-12-5 | 25   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

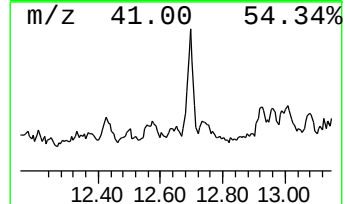
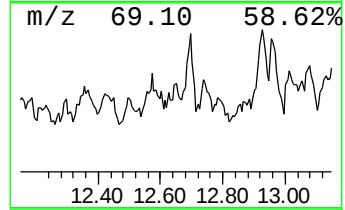
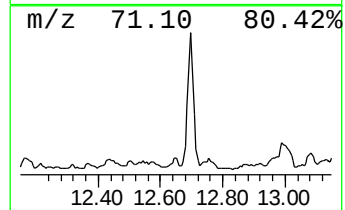
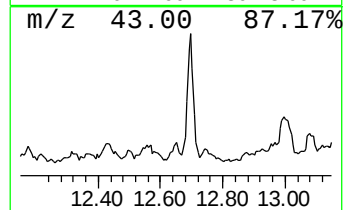
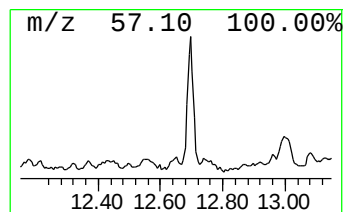
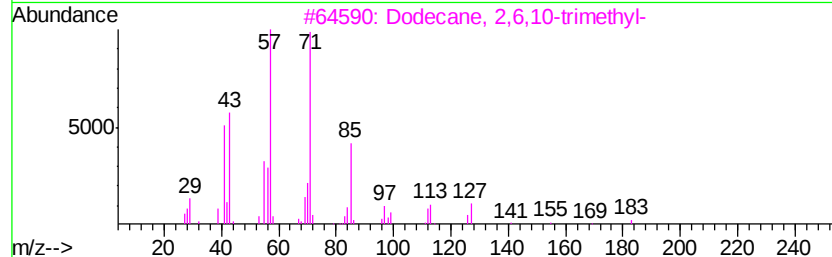
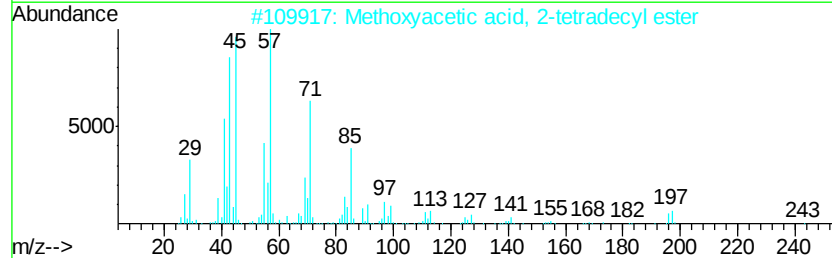
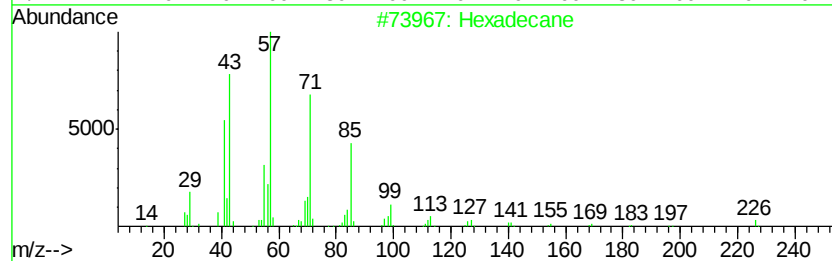
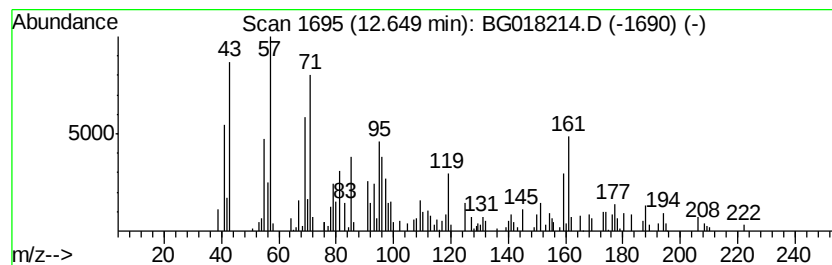
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 40

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.65 | 4.10 ng/ul | 143209 | Acenaphthene-d10 | 14.17 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Hexadecane                          | 226 | C16H34   | 000544-76-3  | 27   |
| 2    |    |   | Methoxyacetic acid, 2-tetradecyl... | 286 | C17H34O3 | 1000282-04-8 | 27   |
| 3    |    |   | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32   | 003891-98-3  | 27   |
| 4    |    |   | Tridecane, 6-propyl-                | 226 | C16H34   | 055045-10-8  | 27   |
| 5    |    |   | Cyclohexanol, 5-methyl-2-(1-meth... | 156 | C10H20O  | 023283-97-8  | 27   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

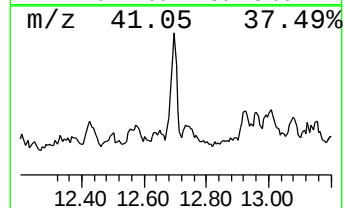
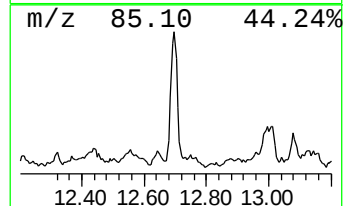
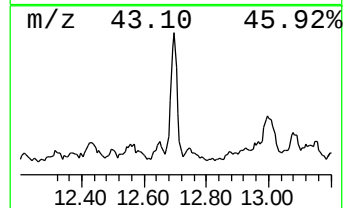
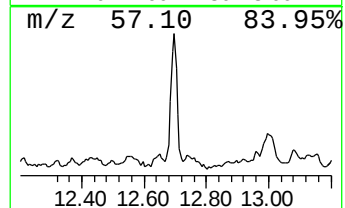
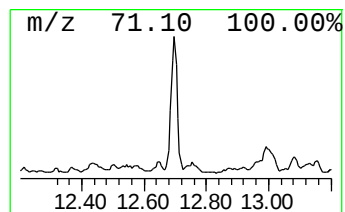
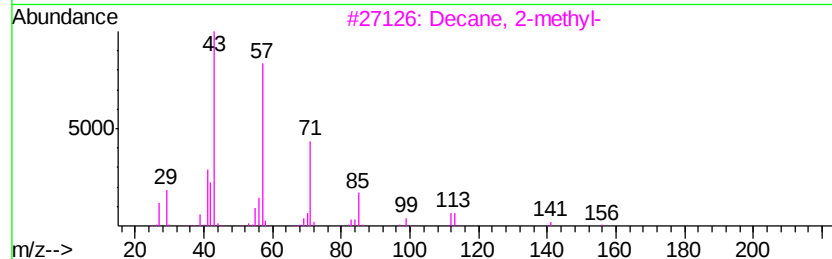
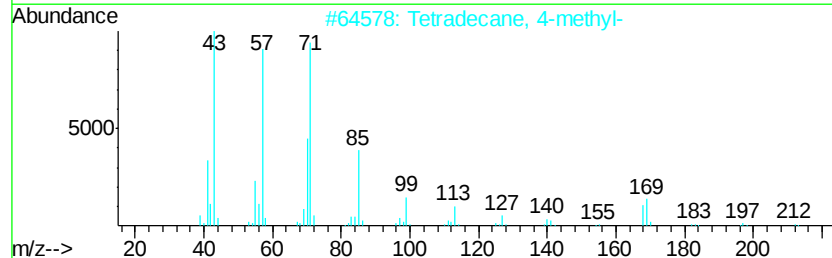
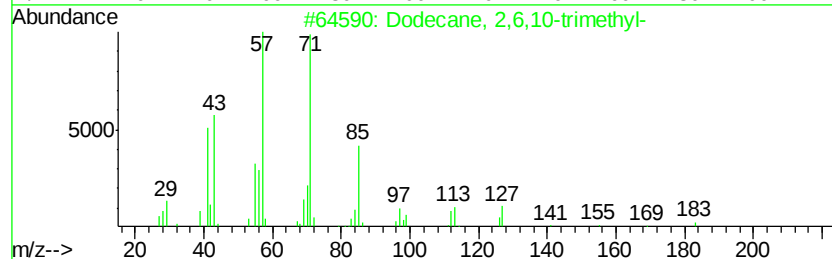
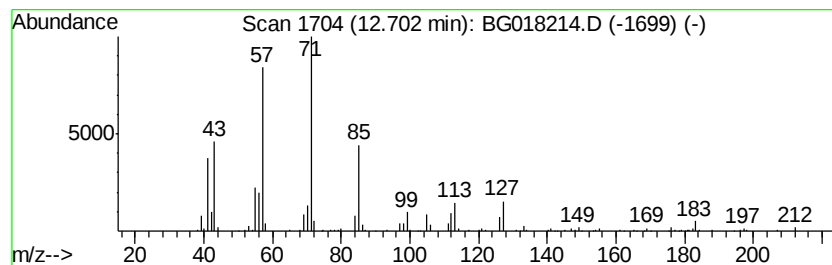
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 12.70 | 15.19 ng/ul | 530283 | Acenaphthene-d10 | 14.17 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 91   |
| 2    |    |   | Tetradecane, 4-methyl-      | 212 | C15H32  | 025117-24-2 | 81   |
| 3    |    |   | Decane, 2-methyl-           | 156 | C11H24  | 006975-98-0 | 72   |
| 4    |    |   | Dodecane, 4-methyl-         | 184 | C13H28  | 006117-97-1 | 68   |
| 5    |    |   | Dodecane, 4,6-dimethyl-     | 198 | C14H30  | 061141-72-8 | 68   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

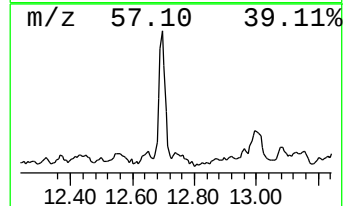
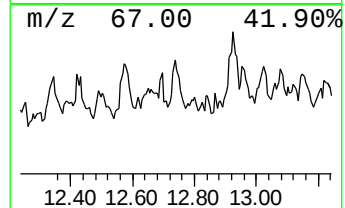
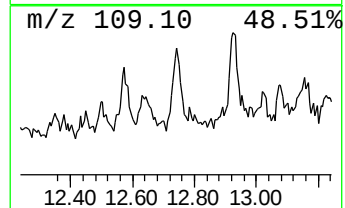
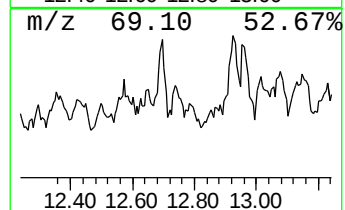
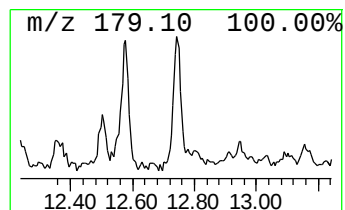
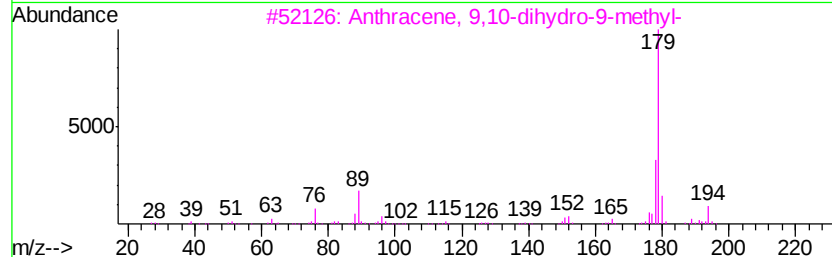
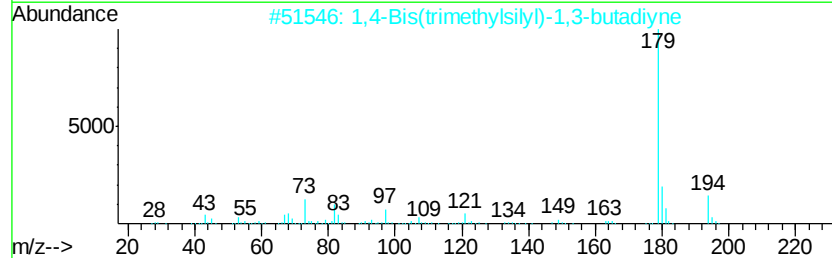
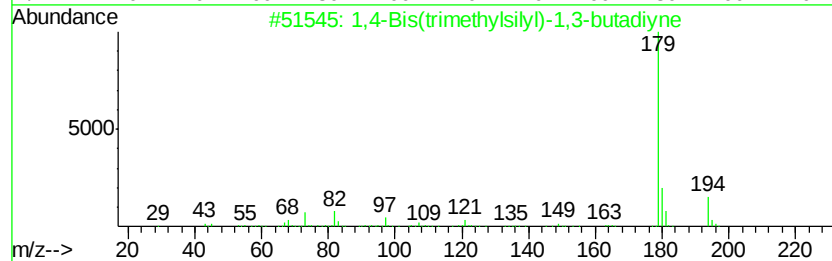
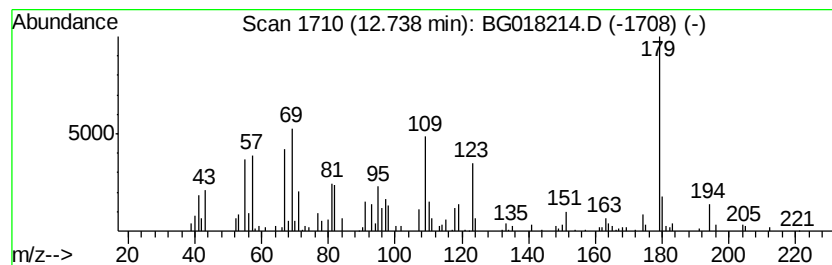
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 unknown-09 Concentration Rank 35

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.74 | 4.59 ng/ul | 160195 | Acenaphthene-d10 | 14.17 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 1,4-Bis(trimethylsilyl)-1,3-buta... | 194 | C10H18Si2    | 004526-07-2 | 27      |      |      |
| 2    | 1,4-Bis(trimethylsilyl)-1,3-buta... | 194 | C10H18Si2    | 004526-07-2 | 27      |      |      |
| 3    | Anthracene, 9,10-dihydro-9-methyl-  | 194 | C15H14       | 017239-99-5 | 27      |      |      |
| 4    | (p-Butyrylphenoxy)acetic acid       | 222 | C12H14O4     | 001141-96-4 | 16      |      |      |
| 5    | Benzoic acid, 2-acetyl-3-methoxy-   | 194 | C10H10O4     | 120714-42-3 | 12      |      |      |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

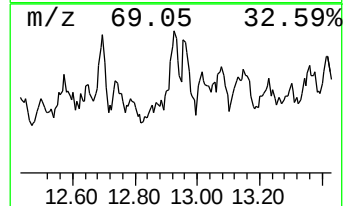
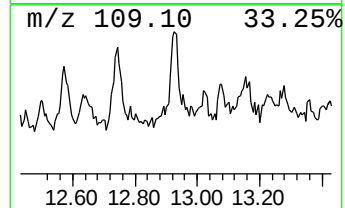
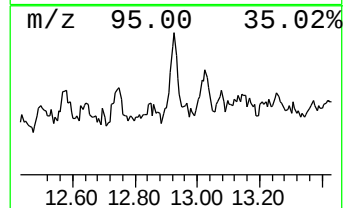
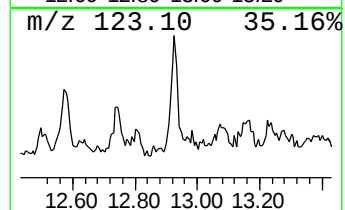
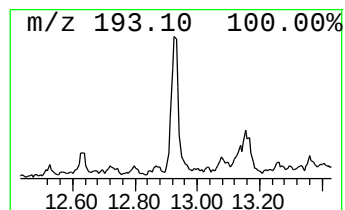
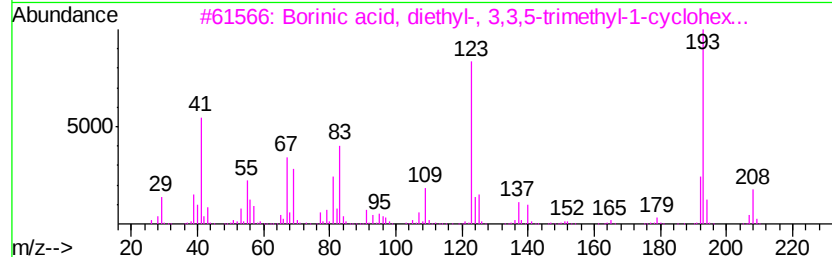
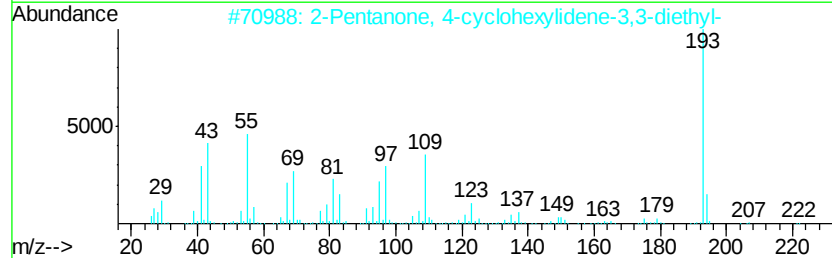
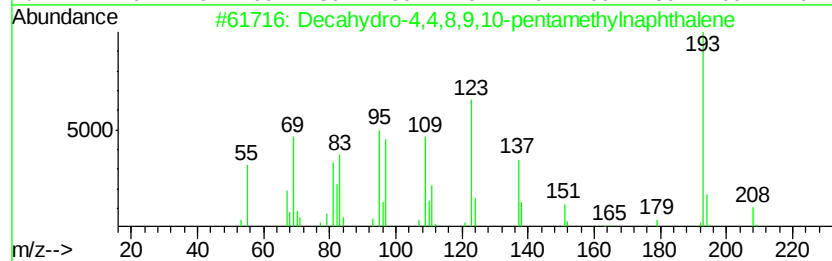
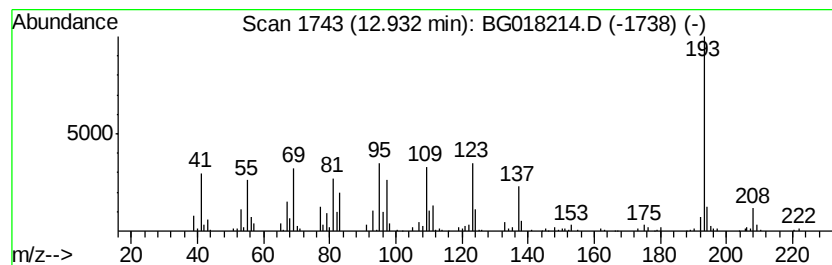
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 Decahydro-4,4,8,9,10-pentam... Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.93 | 6.15 ng/ul | 214702 | Acenaphthene-d10 | 14.17 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Decahydro-4,4,8,9,10-pentamethyl... | 208 | C15H28    | 080655-44-3  | 70   |
| 2    |    | 2-Pentanone, 4-cyclohexylidene-3... | 222 | C15H26O   | 313253-65-5  | 38   |
| 3    |    | Borinic acid, diethyl-, 3,3,5-tr... | 208 | C13H25BO  | 057387-76-5  | 25   |
| 4    |    | 2-Anthracenamine                    | 193 | C14H11N   | 000613-13-8  | 25   |
| 5    |    | 1-(P-FLUOROPHENYL)-4-PIPERIDONE     | 193 | C11H12FNO | 1000238-56-7 | 25   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

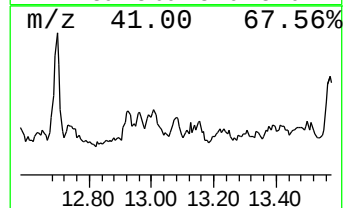
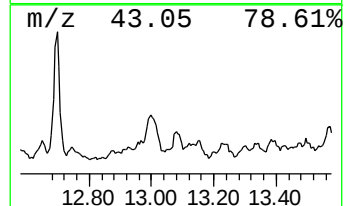
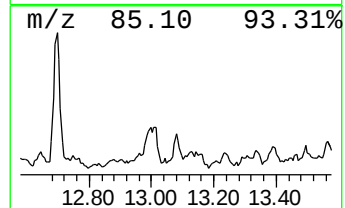
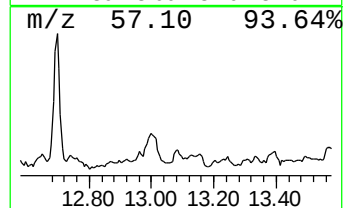
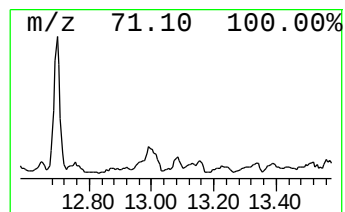
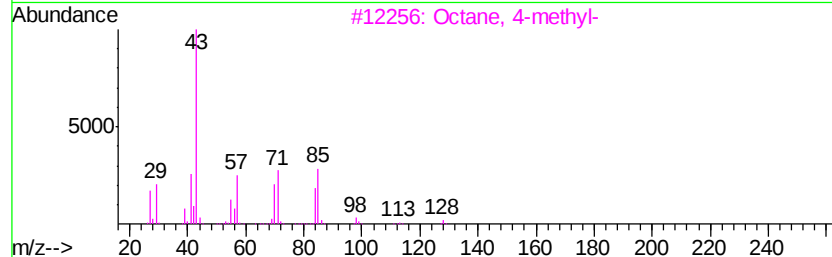
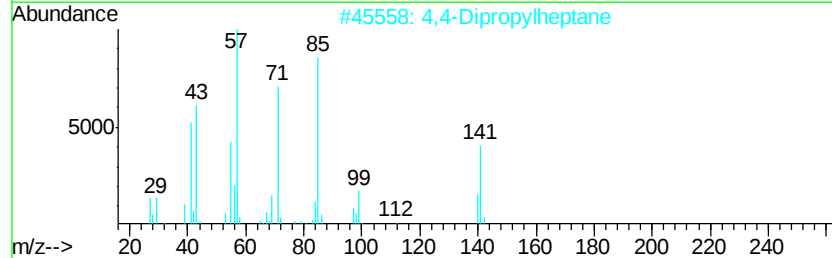
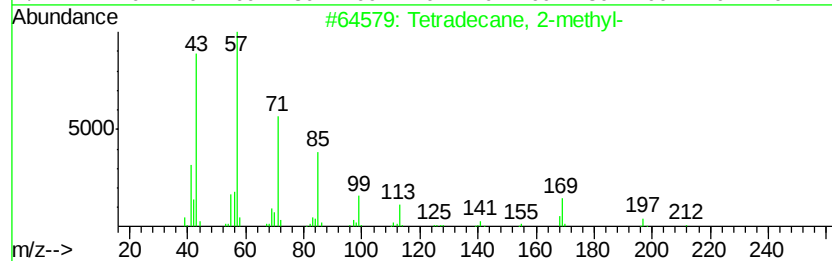
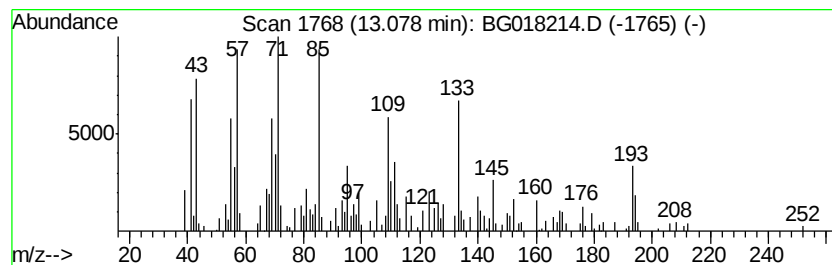
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 43

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.08 | 3.57 ng/ul | 124732 | Acenaphthene-d10 | 14.17 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane, 2-methyl- | 212 | C15H32  | 001560-95-8 | 46   |
| 2    |    |   | 4,4-Dipropylheptane    | 184 | C13H28  | 017312-72-0 | 38   |
| 3    |    |   | Octane, 4-methyl-      | 128 | C9H20   | 002216-34-4 | 35   |
| 4    |    |   | Pentadecane            | 212 | C15H32  | 000629-62-9 | 35   |
| 5    |    |   | Dodecane, 1-iodo-      | 296 | C12H25I | 004292-19-7 | 35   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

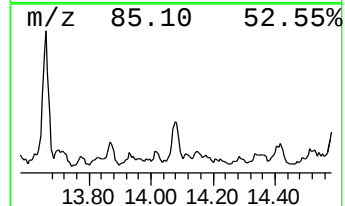
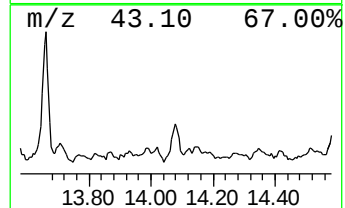
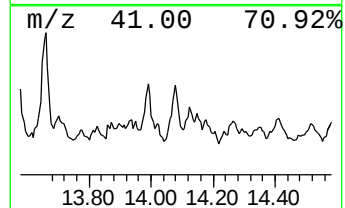
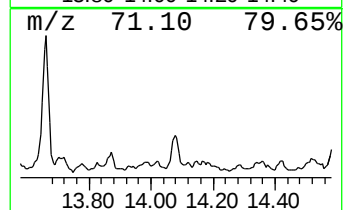
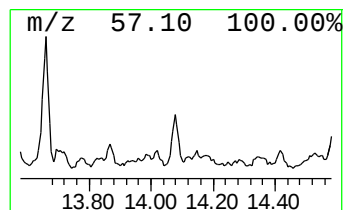
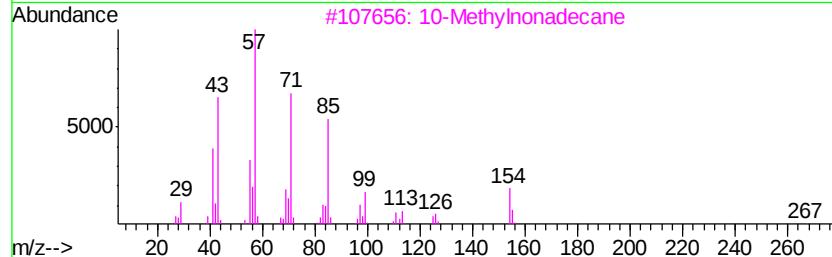
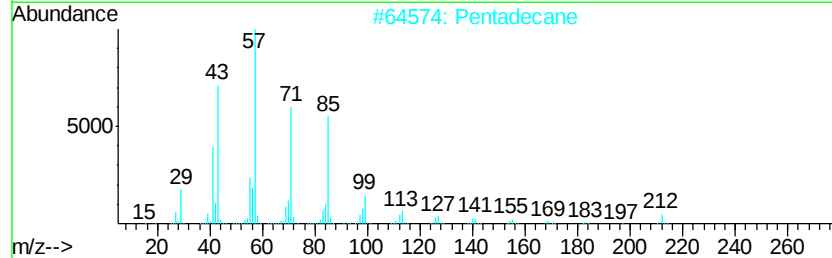
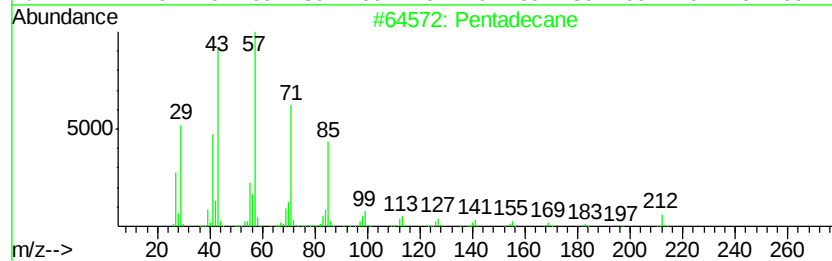
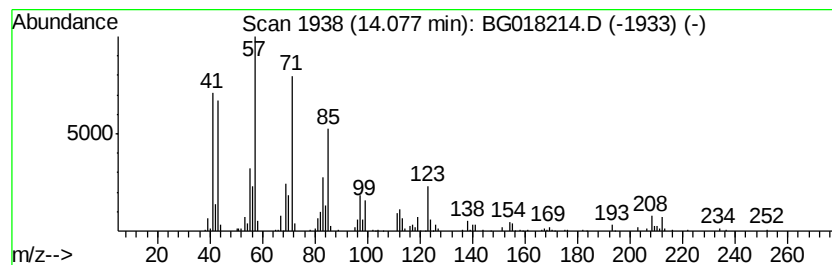
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.08 | 7.25 ng/ul | 253259 | Acenaphthene-d10 | 14.17 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane         | 212 | C15H32  | 000629-62-9 | 89   |
| 2    |    | Pentadecane         | 212 | C15H32  | 000629-62-9 | 89   |
| 3    |    | 10-Methylnonadecane | 282 | C20H42  | 056862-62-5 | 72   |
| 4    |    | Tritetracontane     | 605 | C43H88  | 007098-21-7 | 72   |
| 5    |    | Pentadecane         | 212 | C15H32  | 000629-62-9 | 70   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

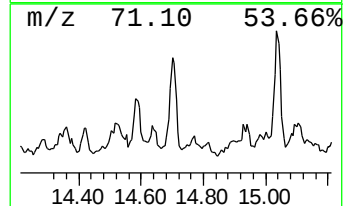
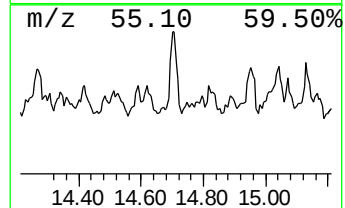
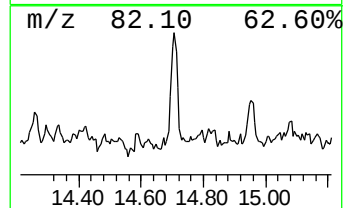
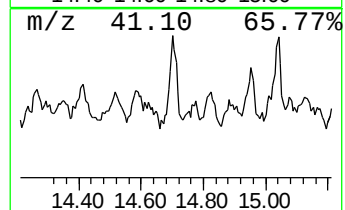
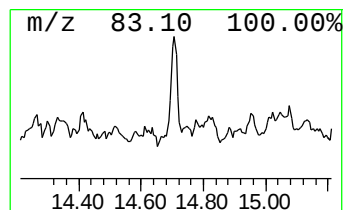
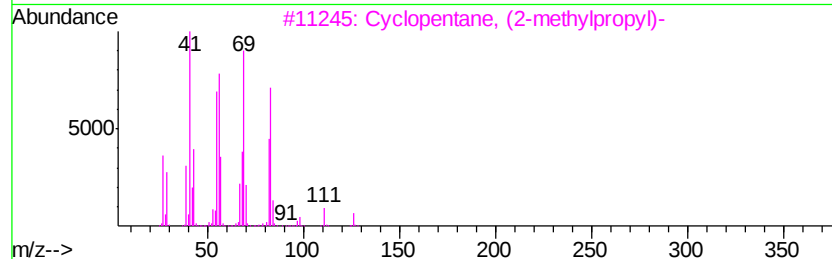
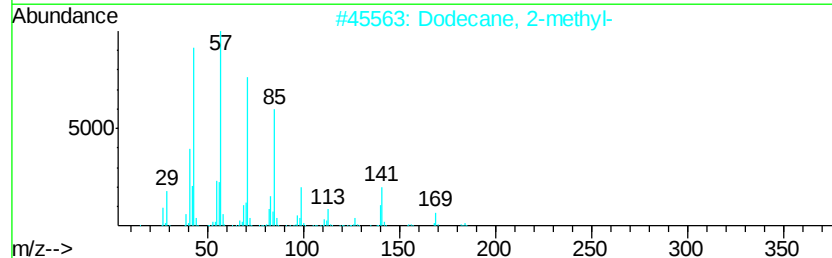
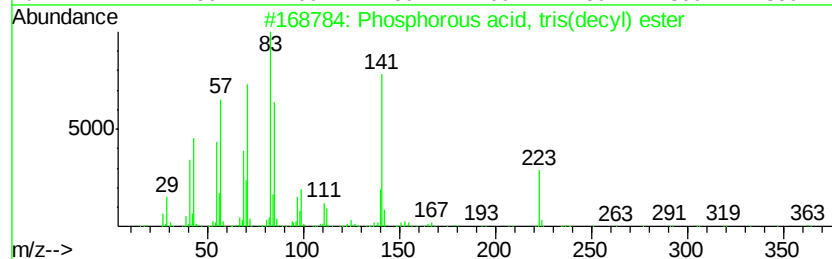
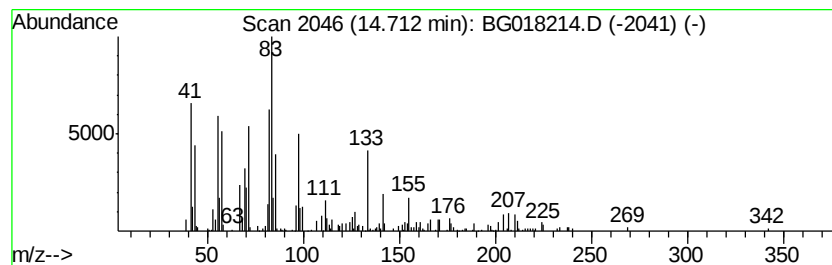
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 unknown-11 Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.71 | 9.29 ng/ul | 324420 | Acenaphthene-d10 | 14.17 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Phosphorous acid, tris(decyl) ester | 502 | C30H63O3P | 002929-86-4 | 49   |
| 2    |    | Dodecane, 2-methyl-                 | 184 | C13H28    | 001560-97-0 | 42   |
| 3    |    | Cyclopentane, (2-methylpropyl)-     | 126 | C9H18     | 003788-32-7 | 38   |
| 4    |    | Diisoomylene                        | 140 | C10H20    | 054063-09-1 | 38   |
| 5    |    | Bemegride                           | 155 | C8H13NO2  | 000064-65-3 | 30   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

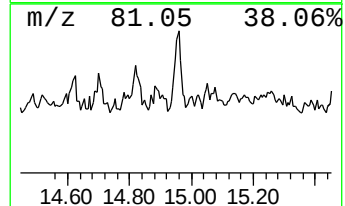
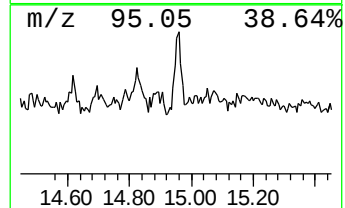
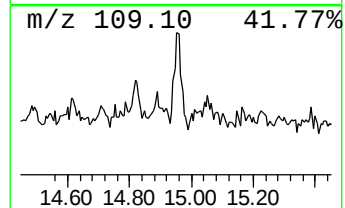
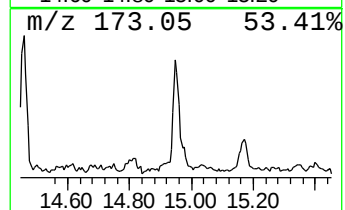
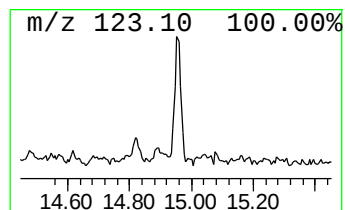
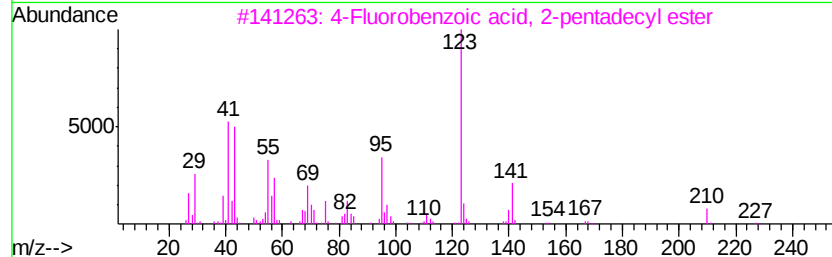
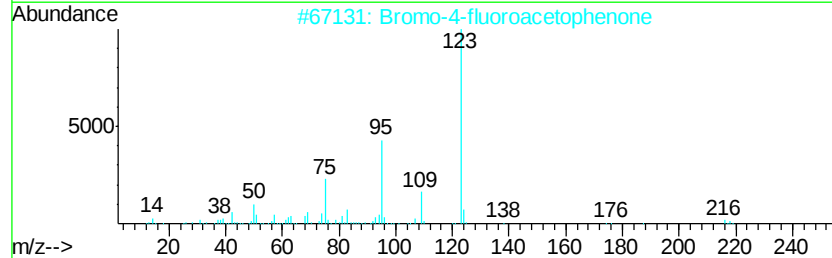
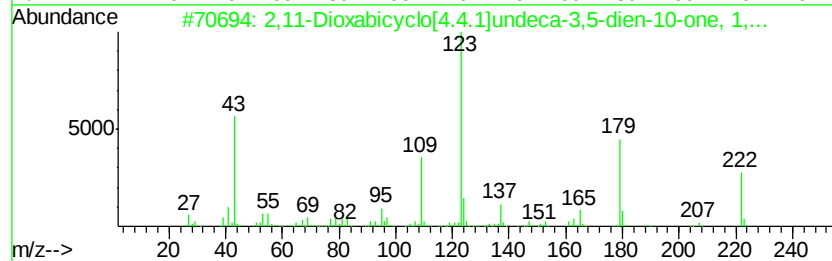
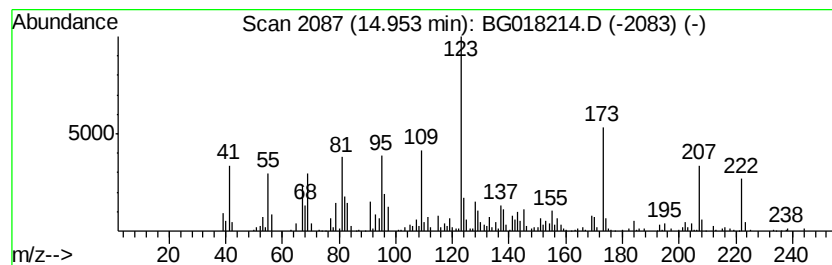
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 unknown-12 Concentration Rank 22

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.95 | 8.09 ng/ul | 282509 | Acenaphthene-d10 | 14.17 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2,11-Dioxabicyclo[4.4.1]undeca-3... | 222 | C13H18O3     | 070412-52-1  | 49      |      |      |
| 2    | Bromo-4-fluoroacetophenone          | 216 | C8H6BrFO     | 000403-29-2  | 38      |      |      |
| 3    | 4-Fluorobenzoic acid, 2-pentadec... | 350 | C22H35FO2    | 1000280-68-3 | 35      |      |      |
| 4    | 3-Fluorobenzoic acid, 2-tridecyl... | 322 | C20H31FO2    | 1000280-60-2 | 35      |      |      |
| 5    | .beta.-Chloro-para-fluoropropiop... | 186 | C9H8ClFO     | 000347-93-3  | 35      |      |      |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

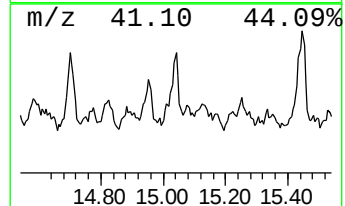
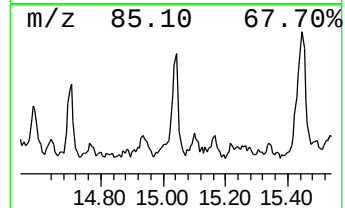
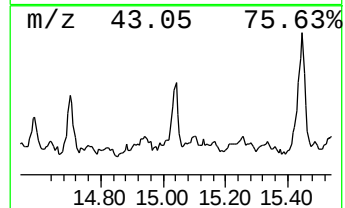
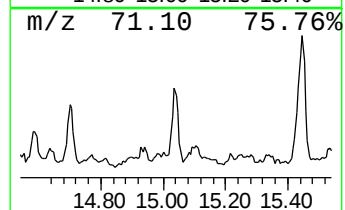
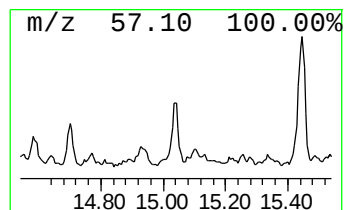
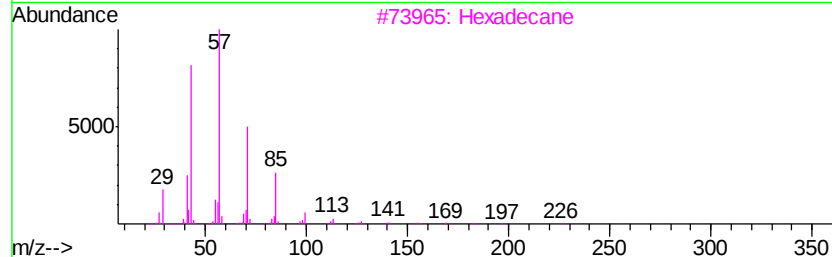
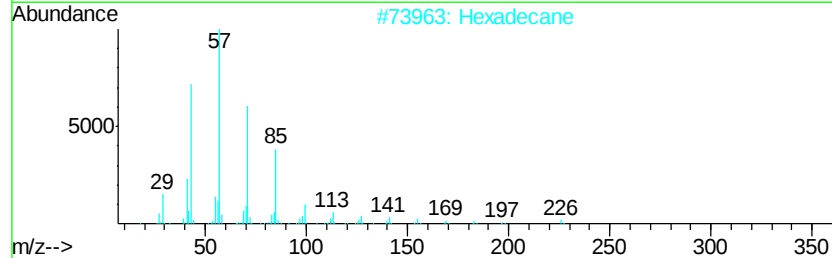
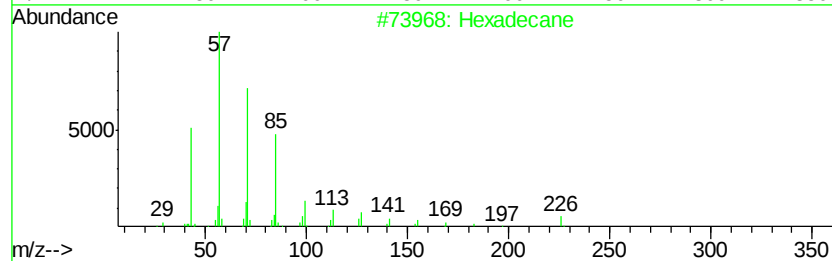
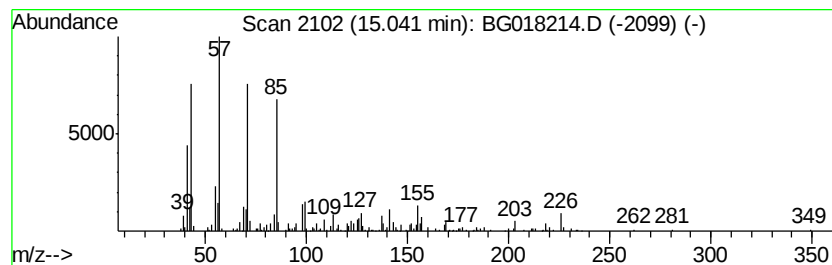
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.04 | 4.82 ng/ul | 168341 | Acenaphthene-d10 | 14.17 |

| Hit# | of          | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-------------|---|--------------|-----|---------|-------------|------|
| 1    | Hexadecane  |   |              | 226 | C16H34  | 000544-76-3 | 93   |
| 2    | Hexadecane  |   |              | 226 | C16H34  | 000544-76-3 | 91   |
| 3    | Hexadecane  |   |              | 226 | C16H34  | 000544-76-3 | 90   |
| 4    | Hexadecane  |   |              | 226 | C16H34  | 000544-76-3 | 83   |
| 5    | Pentadecane |   |              | 212 | C15H32  | 000629-62-9 | 81   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

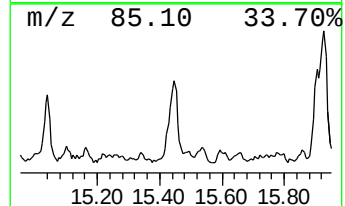
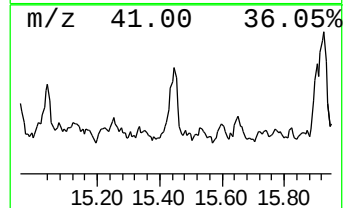
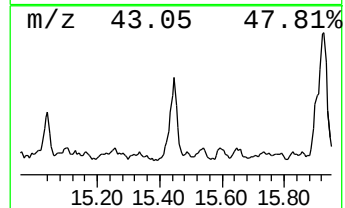
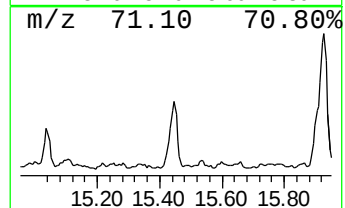
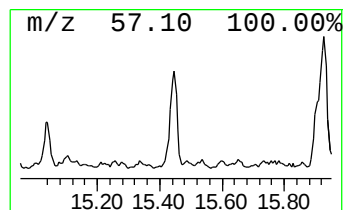
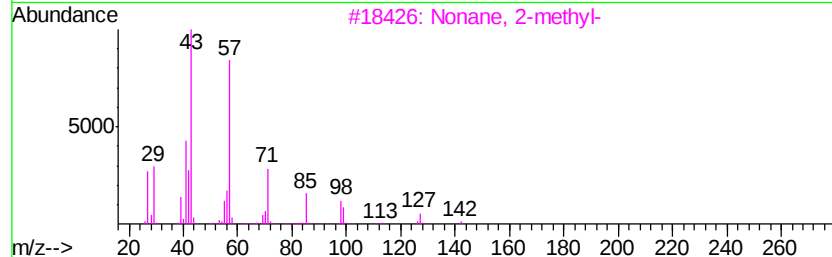
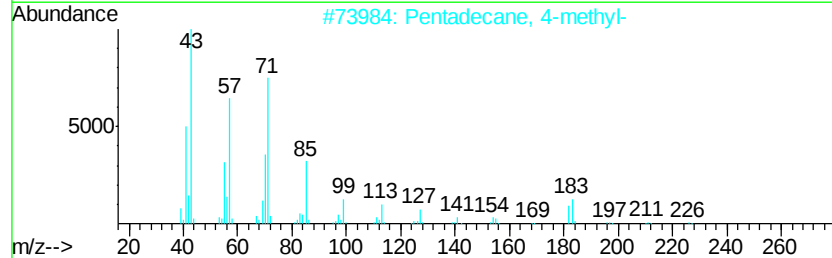
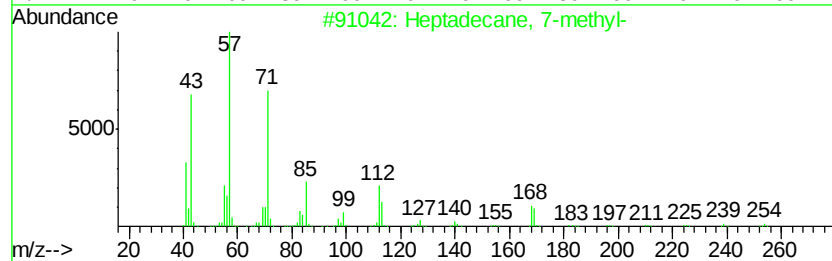
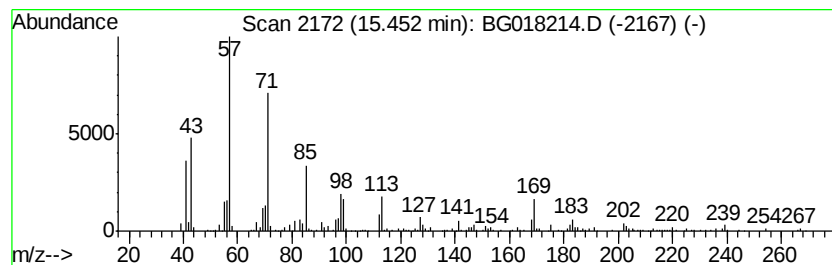
Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.    | EstConc     | Area                   | Relative to ISTD | R.T.           |
|---------|-------------|------------------------|------------------|----------------|
| 15.45   | 13.32 ng/ul | 464960                 | Acenaphthene-d10 | 14.17          |
| Hit# of | 5           | Tentative ID           | MW MolForm       | CAS# Qual      |
| 1       |             | Heptadecane, 7-methyl- | 254 C18H38       | 020959-33-5 83 |
| 2       |             | Pentadecane, 4-methyl- | 226 C16H34       | 002801-87-8 64 |
| 3       |             | Nonane, 2-methyl-      | 142 C10H22       | 000871-83-0 64 |
| 4       |             | Tridecane, 5-propyl-   | 226 C16H34       | 055045-11-9 64 |
| 5       |             | Pentadecane, 7-methyl- | 226 C16H34       | 006165-40-8 64 |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

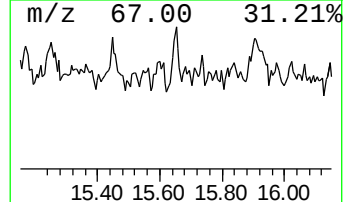
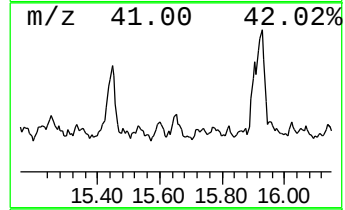
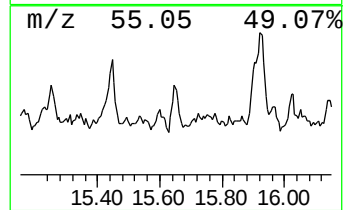
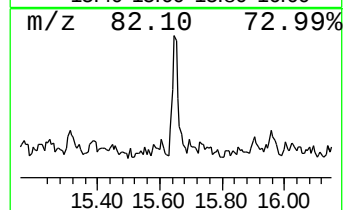
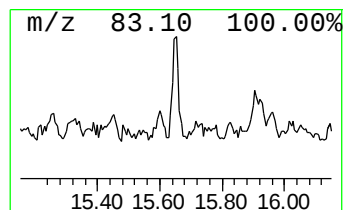
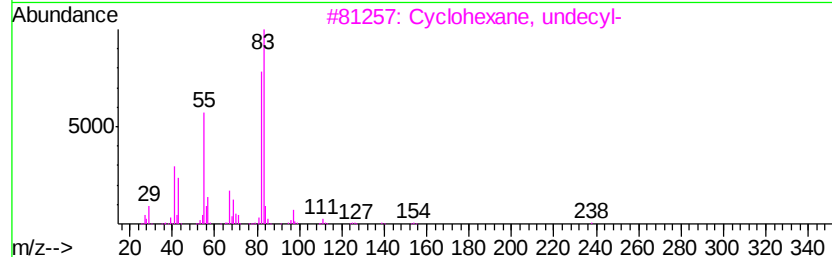
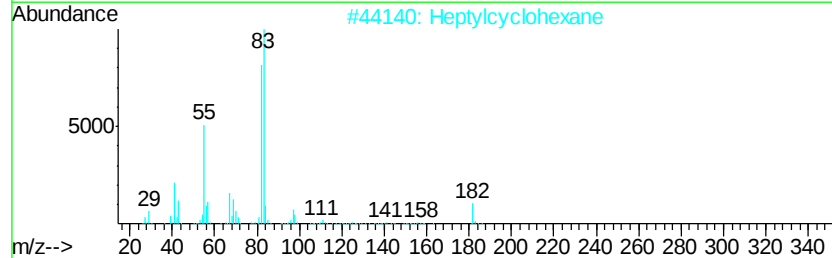
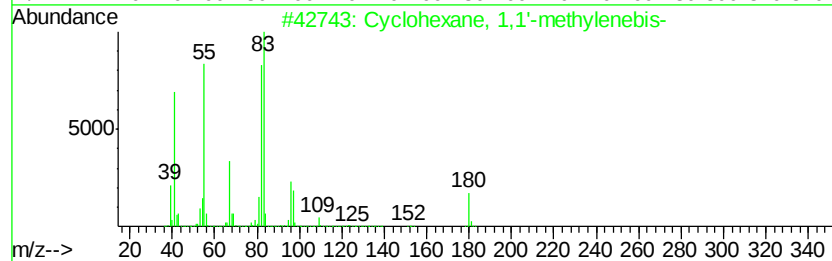
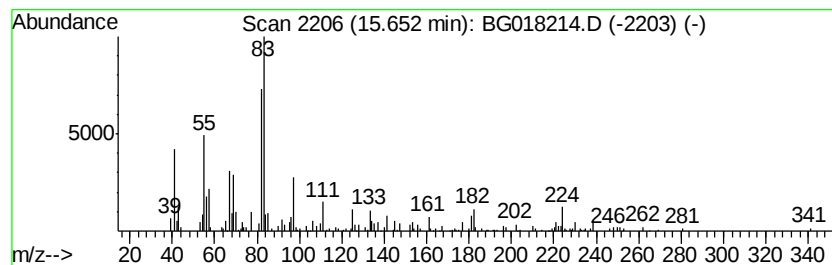
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 37 Cyclohexane, 1,1'-methylene... Concentration Rank 26

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.65 | 6.09 ng/ul | 154240 | Phenanthrene-d10 | 16.92 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,1'-methylenebis- | 180 | C13H24  | 003178-23-2 | 52   |
| 2    |    |   | Heptylcyclohexane               | 182 | C13H26  | 005617-41-4 | 52   |
| 3    |    |   | Cyclohexane, undecyl-           | 238 | C17H34  | 054105-66-7 | 50   |
| 4    |    |   | Cyclohexane, decyl-             | 224 | C16H32  | 001795-16-0 | 50   |
| 5    |    |   | Cyclohexane, octyl-             | 196 | C14H28  | 001795-15-9 | 49   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

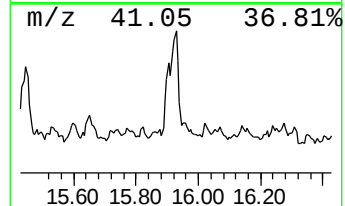
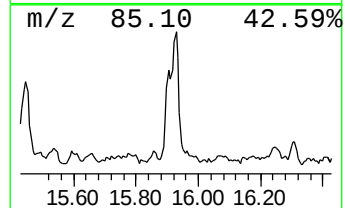
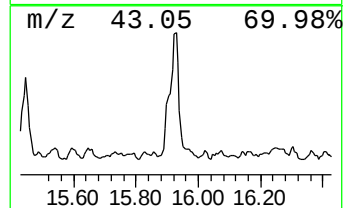
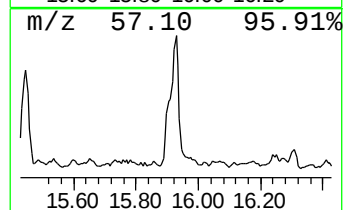
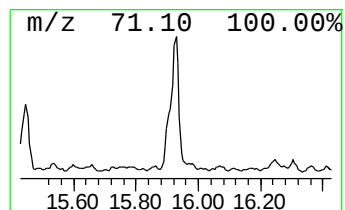
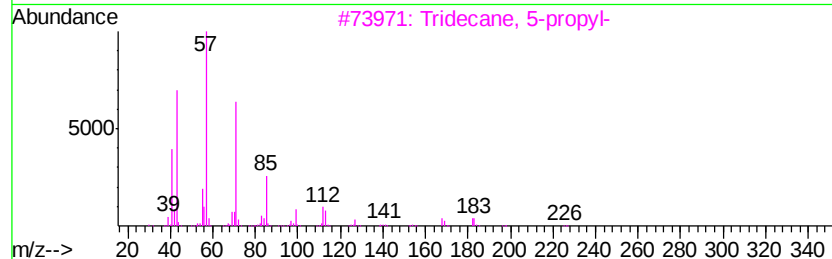
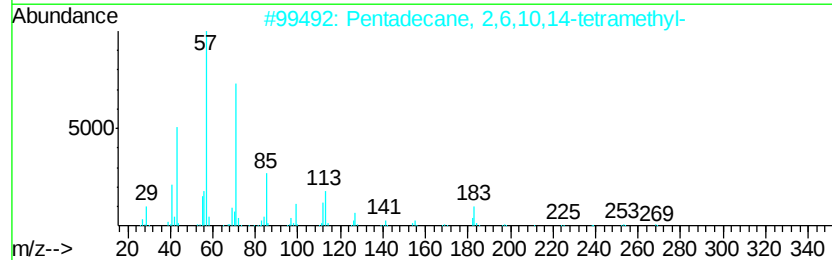
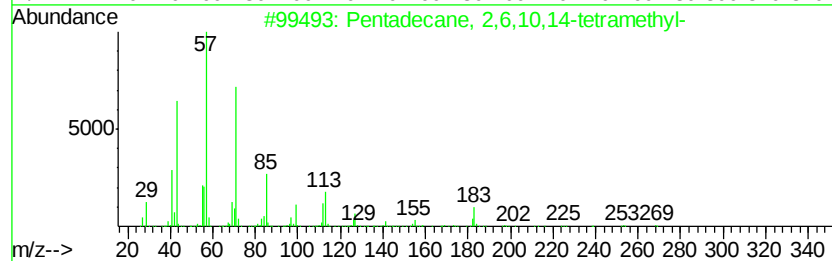
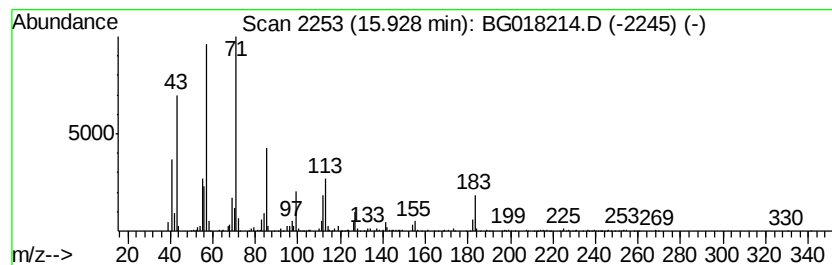
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.93 | 42.76 ng/ul | 1083290 | Phenanthrene-d10 | 16.92 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6  | 95   |
| 2    |    | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6  | 93   |
| 3    |    | Tridecane, 5-propyl-                | 226 | C16H34  | 055045-11-9  | 81   |
| 4    |    | 7-Methyl-octadecane                 | 268 | C19H40  | 1000192-63-3 | 78   |
| 5    |    | Tridecane, 6-propyl-                | 226 | C16H34  | 055045-10-8  | 60   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

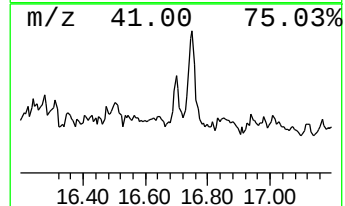
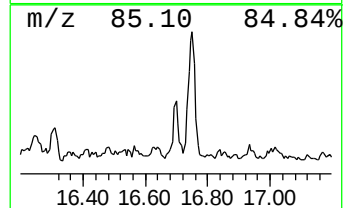
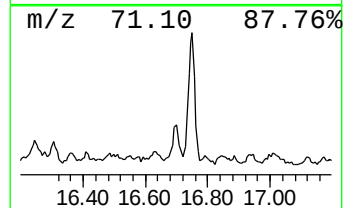
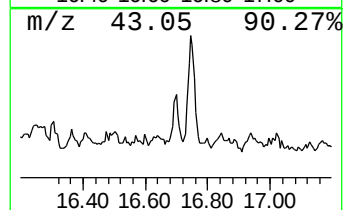
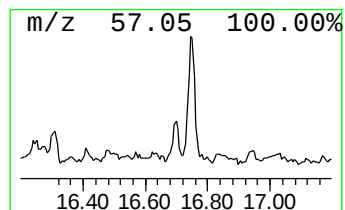
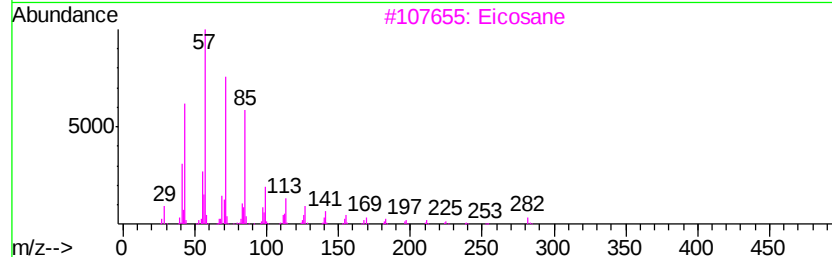
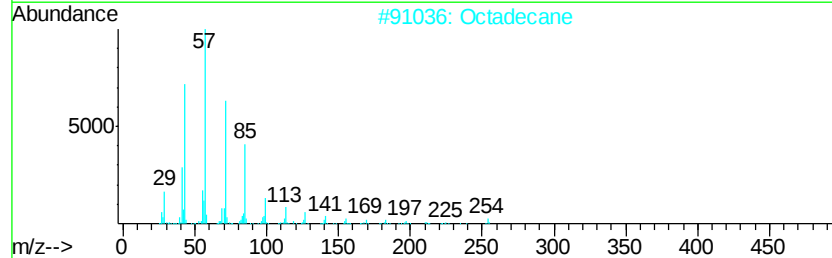
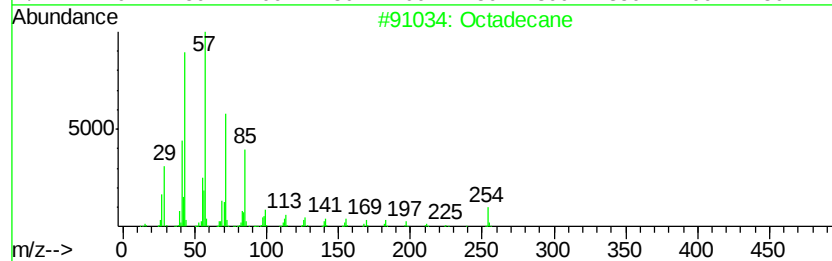
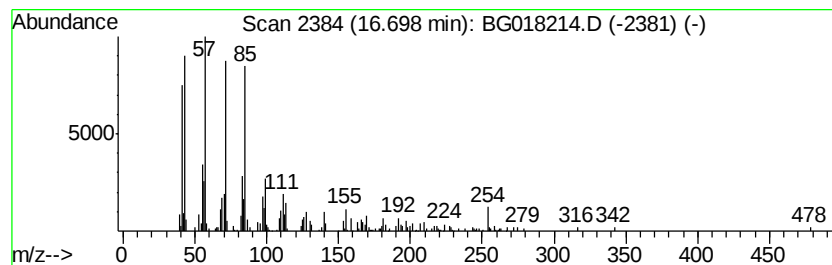
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 39

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.70 | 4.26 ng/ul | 107870 | Phenanthrene-d10 | 16.92 |

| Hit# | of | 5 | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------|-----|---------|-------------|------|
| 1    |    |   | Octadecane       | 254 | C18H38  | 000593-45-3 | 90   |
| 2    |    |   | Octadecane       | 254 | C18H38  | 000593-45-3 | 80   |
| 3    |    |   | Eicosane         | 282 | C20H42  | 000112-95-8 | 72   |
| 4    |    |   | Heneicosane      | 296 | C21H44  | 000629-94-7 | 72   |
| 5    |    |   | Tetratriacontane | 479 | C34H70  | 014167-59-0 | 72   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

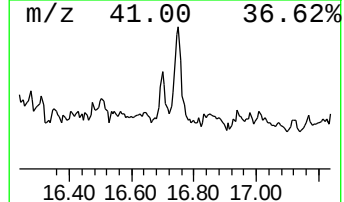
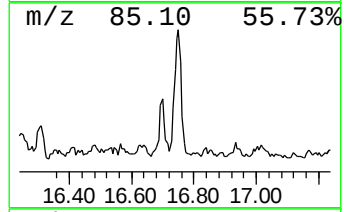
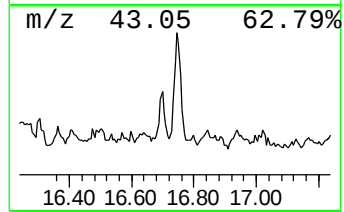
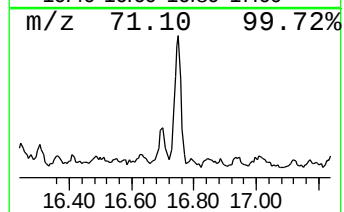
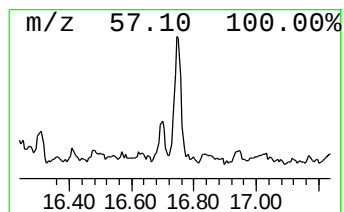
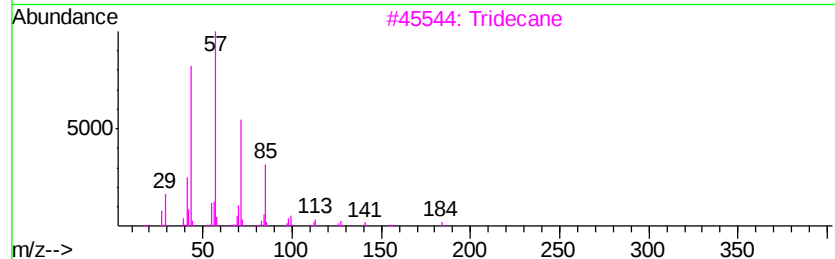
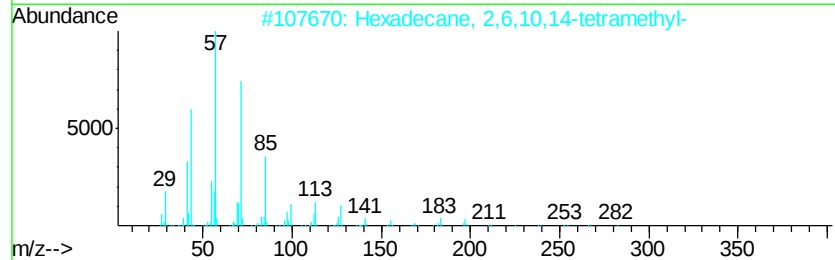
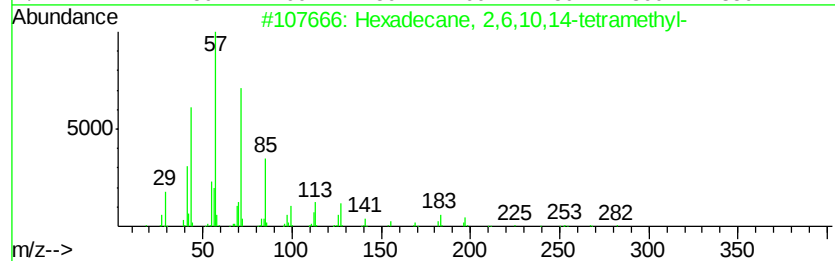
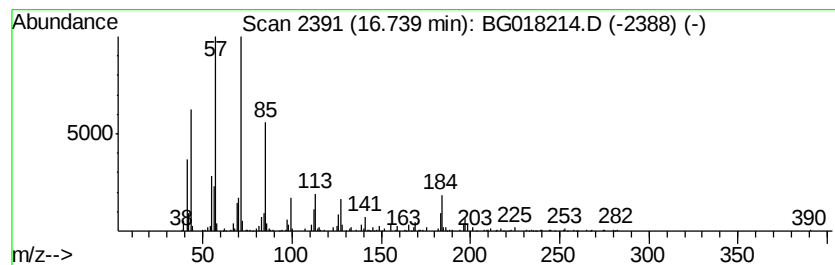
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 16.74 | 19.29 ng/ul | 488722 | Phenanthrene-d10 | 16.92 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 96   |
| 2    |    |   | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 87   |
| 3    |    |   | Tridecane                          | 184 | C13H28  | 000629-50-5 | 87   |
| 4    |    |   | Hexadecane                         | 226 | C16H34  | 000544-76-3 | 87   |
| 5    |    |   | Hexadecane, 2-methyl-              | 240 | C17H36  | 001560-92-5 | 86   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

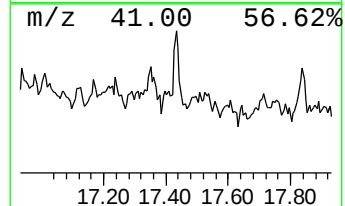
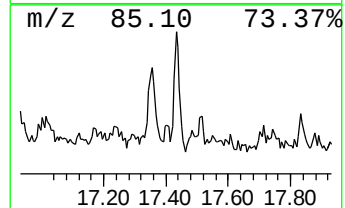
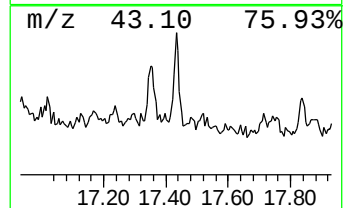
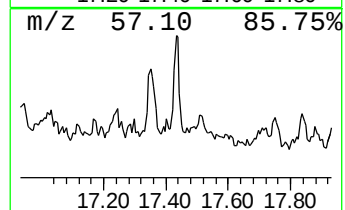
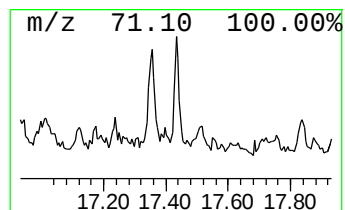
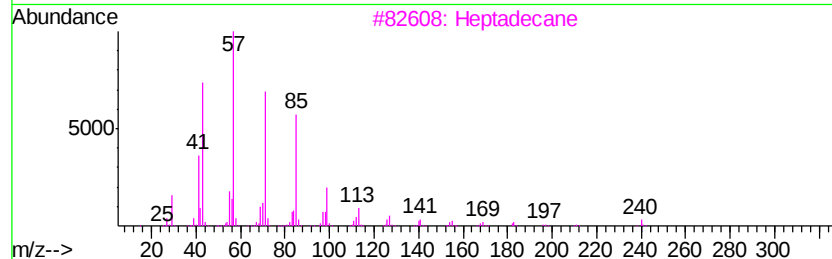
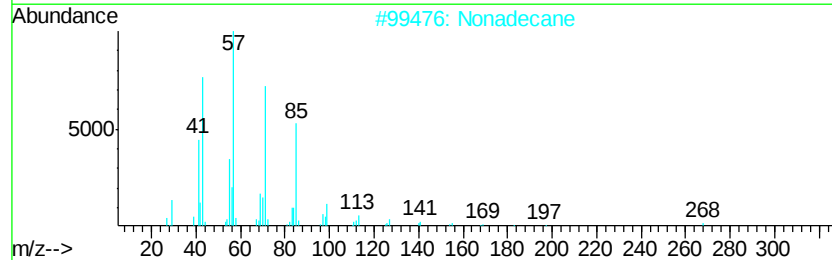
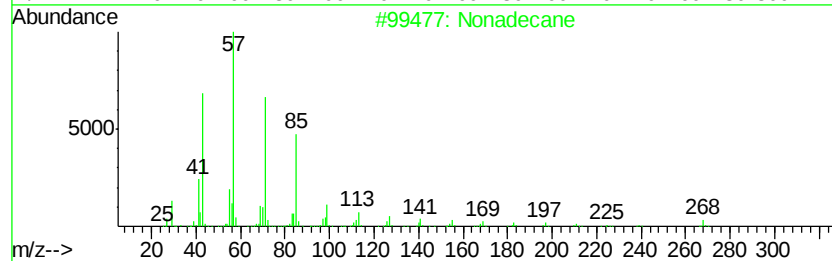
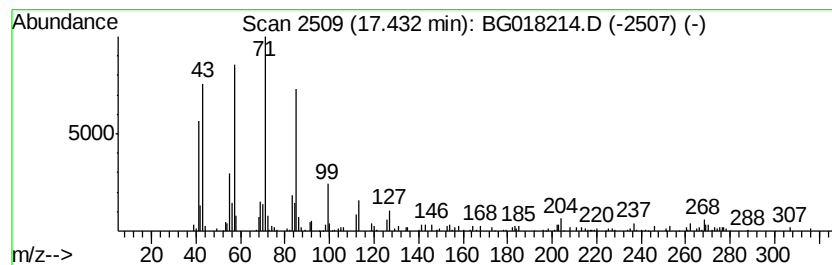
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.43 | 4.58 ng/ul | 116102 | Phenanthrene-d10 | 16.92 |

| Hit# | of | 5 | Tentative ID     | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane       | 268 | C19H40  | 000629-92-5 | 90   |
| 2    |    |   | Nonadecane       | 268 | C19H40  | 000629-92-5 | 87   |
| 3    |    |   | Heptadecane      | 240 | C17H36  | 000629-78-7 | 80   |
| 4    |    |   | Tetratriacontane | 479 | C34H70  | 014167-59-0 | 72   |
| 5    |    |   | Hexacosane       | 366 | C26H54  | 000630-01-3 | 72   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

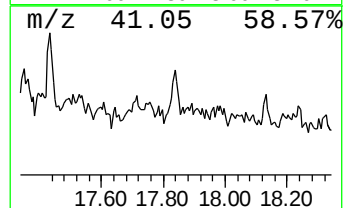
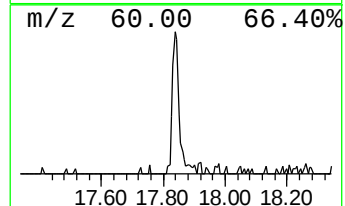
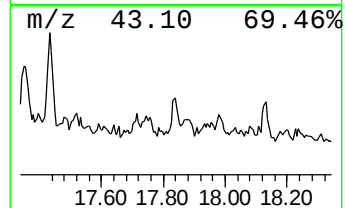
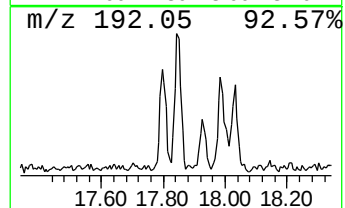
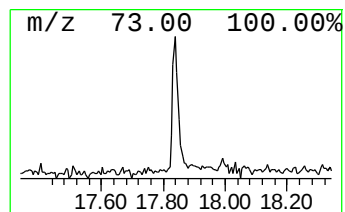
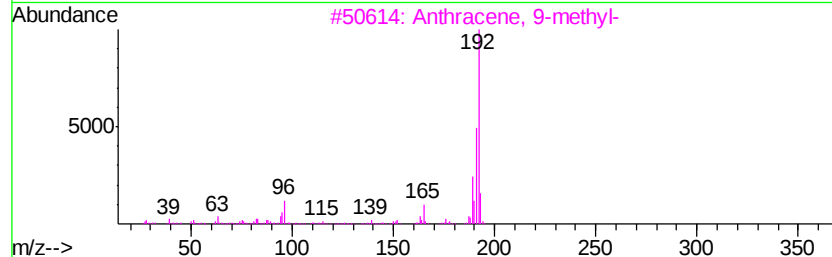
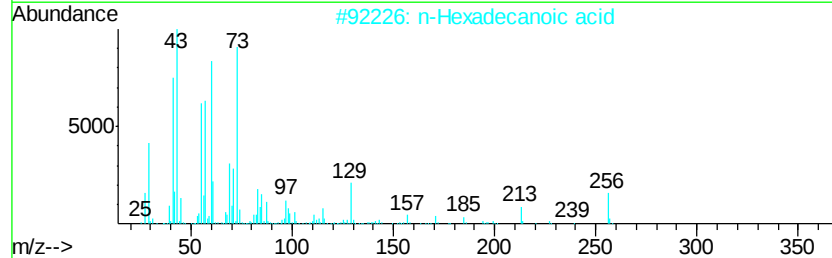
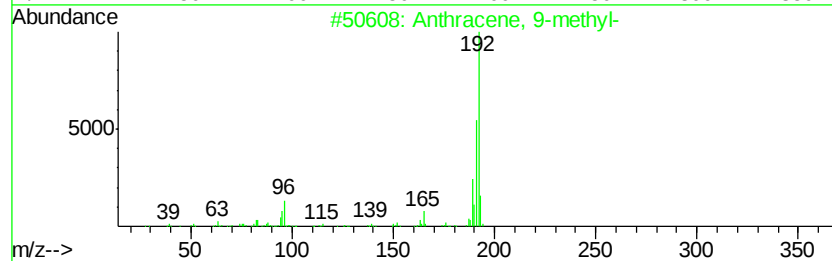
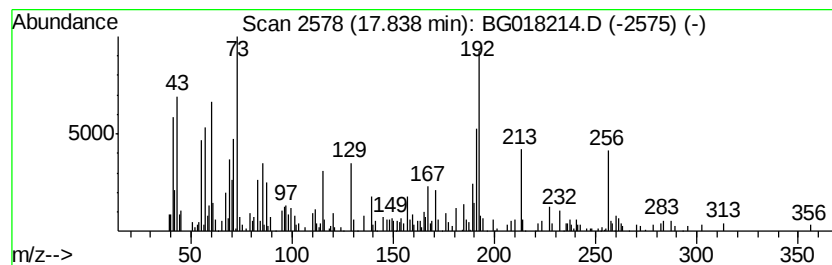
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 42 unknown-13 Concentration Rank 32

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.84 | 4.99 ng/ul | 126470 | Phenanthrene-d10 | 16.92 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------|-----|----------|-------------|------|
| 1    |    |   | Anthracene, 9-methyl-   | 192 | C15H12   | 000779-02-2 | 38   |
| 2    |    |   | n-Hexadecanoic acid     | 256 | C16H32O2 | 000057-10-3 | 38   |
| 3    |    |   | Anthracene, 9-methyl-   | 192 | C15H12   | 000779-02-2 | 38   |
| 4    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12   | 002531-84-2 | 30   |
| 5    |    |   | Anthracene, 1-methyl-   | 192 | C15H12   | 000610-48-0 | 30   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

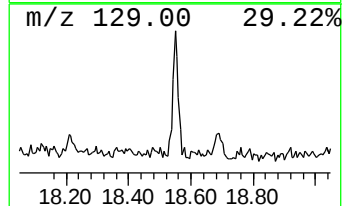
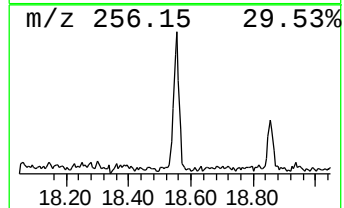
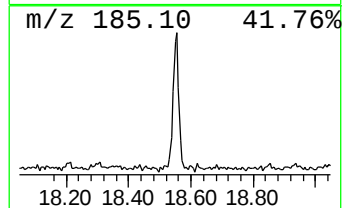
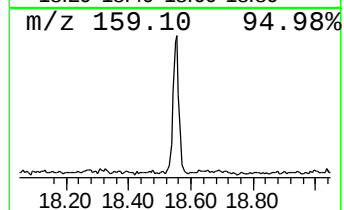
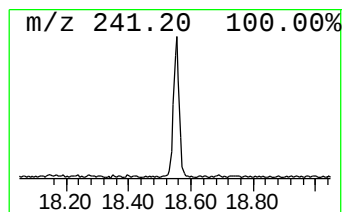
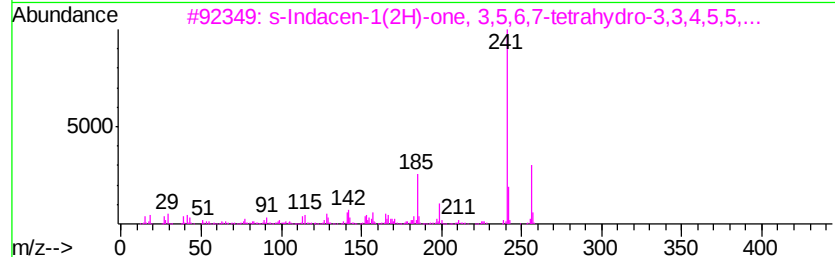
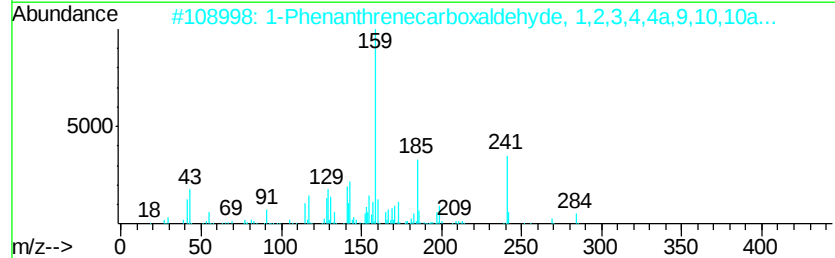
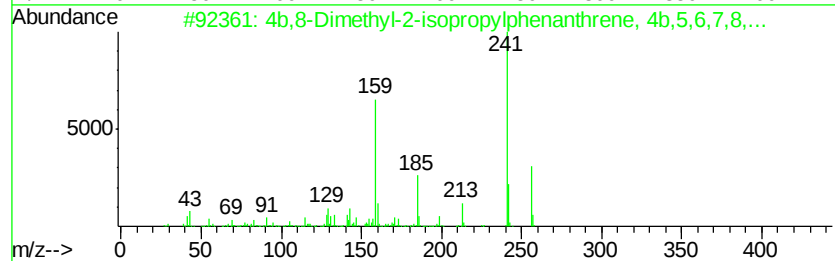
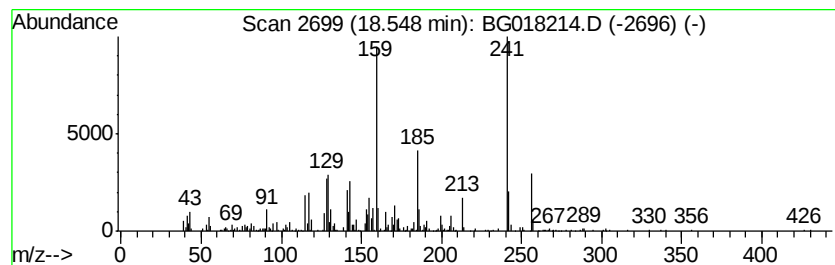
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 43 4b,8-Dimethyl-2-isopropylph... Concentration Rank 15

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 18.55 | 12.27 ng/ul | 310829 | Phenanthrene-d10 | 16.92 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 4b,8-Dimethyl-2-isopropylphenant... | 256 | C19H28  | 1000197-14-1 | 87   |
| 2    |    | 1-Phenanthrenecarboxaldehyde, 1,... | 284 | C20H28O | 024035-50-5  | 83   |
| 3    |    | s-Indacen-1(2H)-one, 3,5,6,7-tet... | 256 | C18H24O | 038754-94-8  | 42   |
| 4    |    | Naphthalene, 6-ethyl-1,2,3,4-tet... | 256 | C19H28  | 301643-35-6  | 38   |
| 5    |    | As-Indacene, 1,2,3,6,7,8-hexahyd... | 256 | C19H28  | 017465-47-3  | 27   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

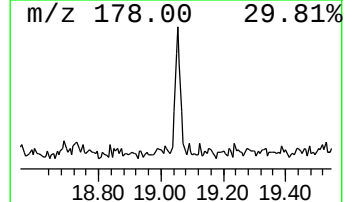
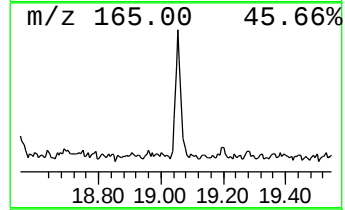
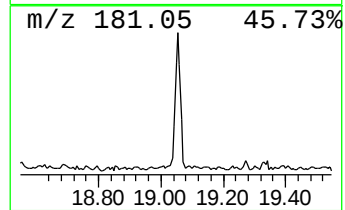
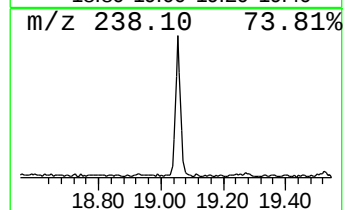
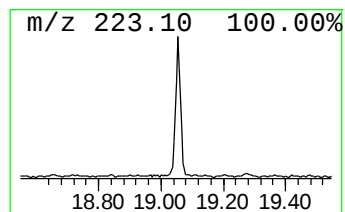
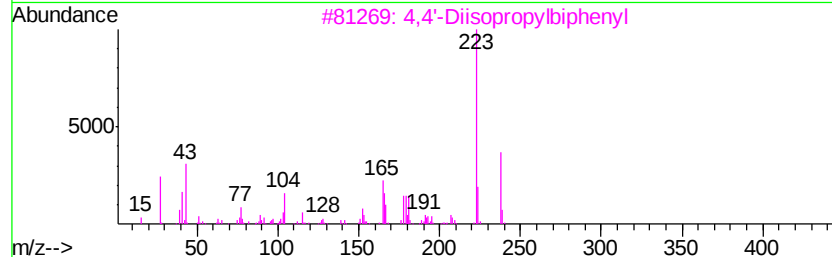
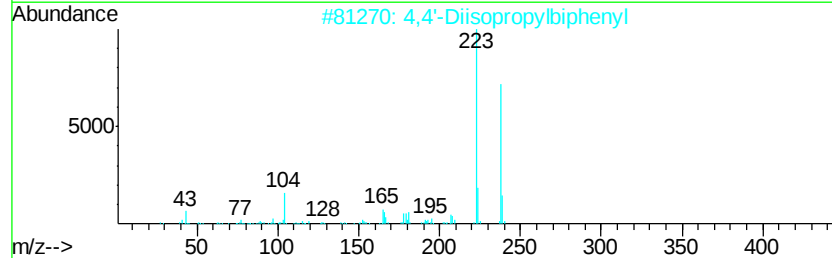
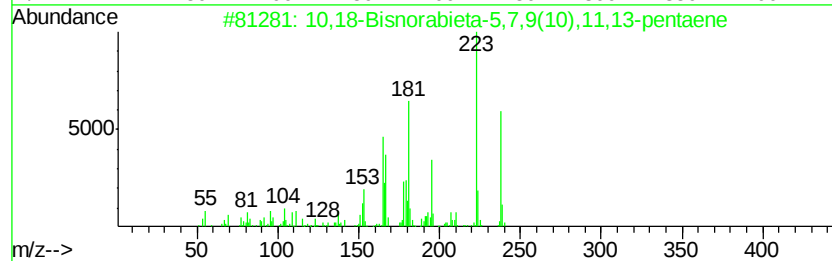
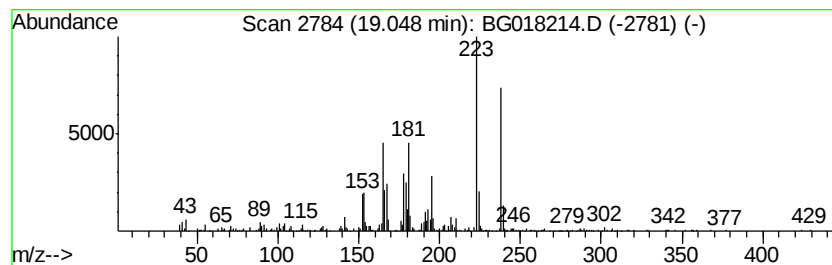
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 44 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 19

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.05 | 8.94 ng/ul | 394643 | Chrysene-d12     | 21.13 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 10,18-Bisnorabieta-5,7,9(10),11,... | 238 | C18H22   | 006566-19-4 | 93   |
| 2    |    |   | 4,4'-Diisopropylbiphenyl            | 238 | C18H22   | 018970-30-4 | 50   |
| 3    |    |   | 4,4'-Diisopropylbiphenyl            | 238 | C18H22   | 018970-30-4 | 50   |
| 4    |    |   | Anthracene, 9,10-dimethoxy-         | 238 | C16H14O2 | 002395-97-3 | 46   |
| 5    |    |   | Anthracene, 1,4-dimethoxy-          | 238 | C16H14O2 | 013076-29-4 | 45   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

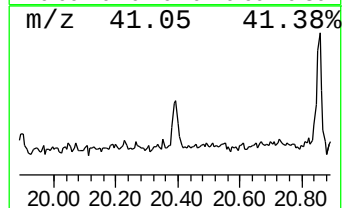
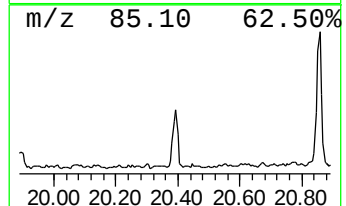
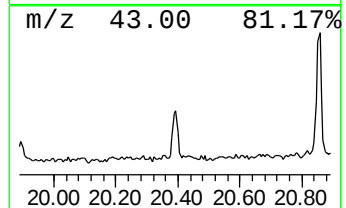
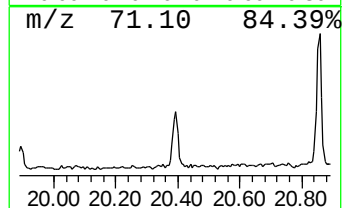
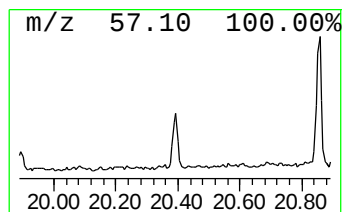
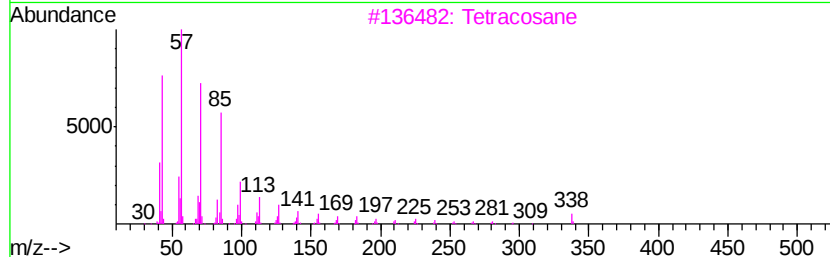
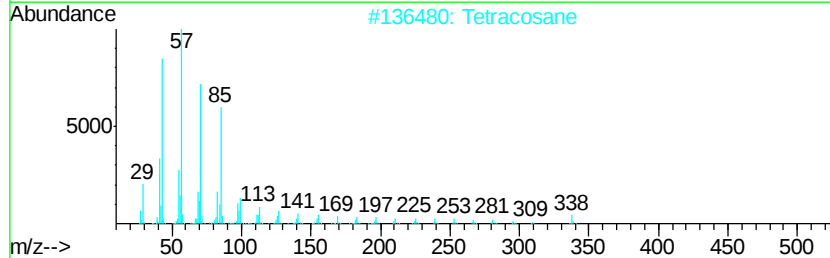
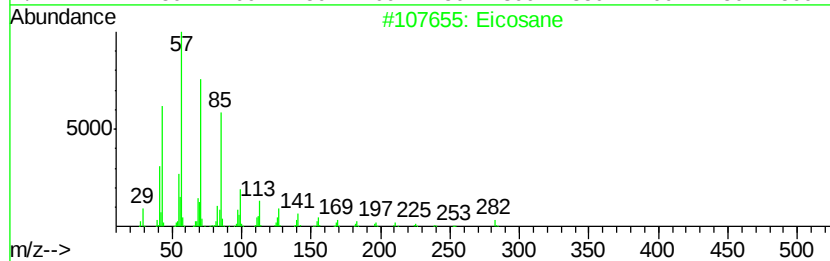
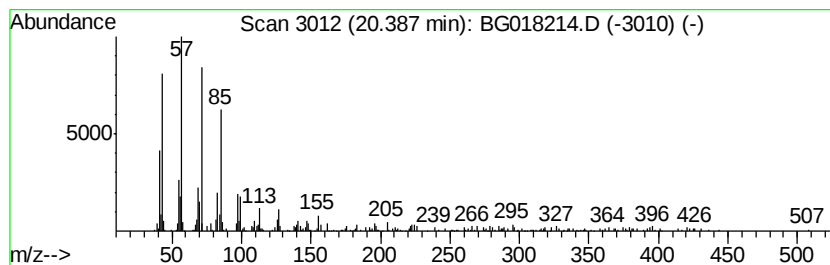
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 44

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.39 | 3.41 ng/ul | 150760 | Chrysene-d12     | 21.13 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane     | 282 | C20H42  | 000112-95-8 | 98   |
| 2    |    | Tetracosane  | 338 | C24H50  | 000646-31-1 | 96   |
| 3    |    | Tetracosane  | 338 | C24H50  | 000646-31-1 | 92   |
| 4    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 91   |
| 5    |    | Tetracosane  | 338 | C24H50  | 000646-31-1 | 87   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

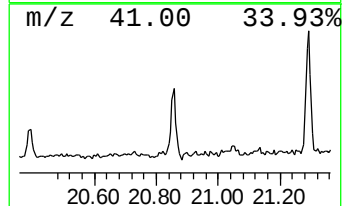
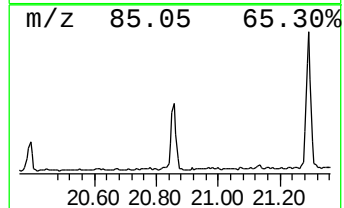
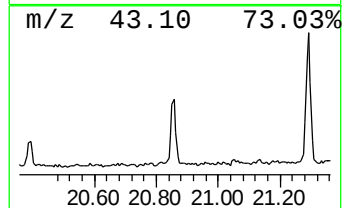
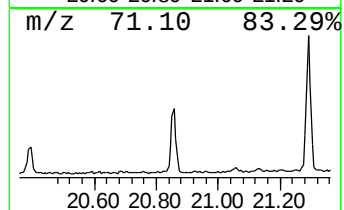
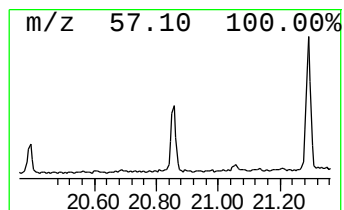
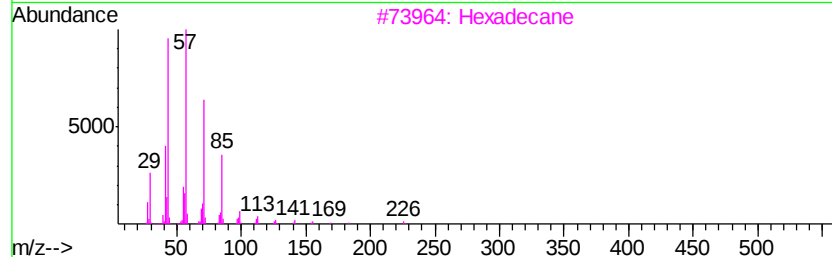
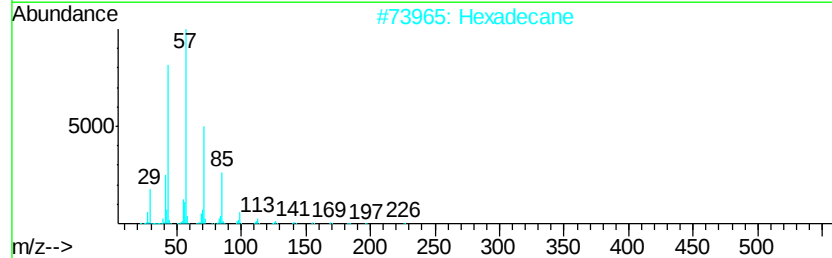
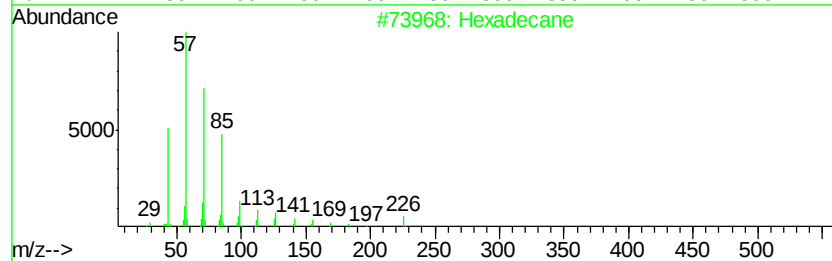
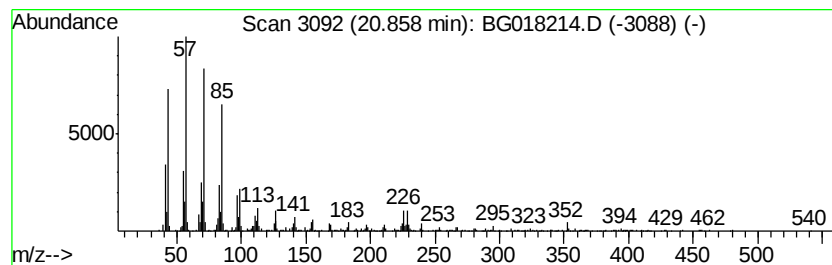
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 20.86 | 10.85 ng/ul | 478938 | Chrysene-d12     | 21.13 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane   | 226 | C16H34  | 000544-76-3 | 97   |
| 2    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 96   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 95   |
| 4    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 95   |
| 5    |    | Octacosane   | 394 | C28H58  | 000630-02-4 | 95   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

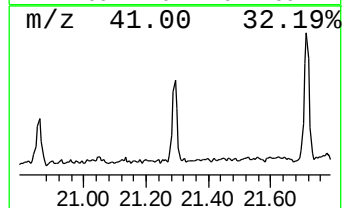
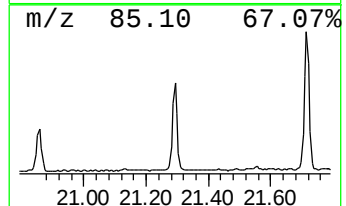
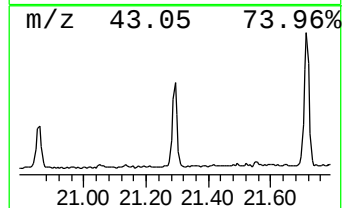
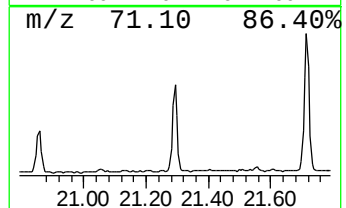
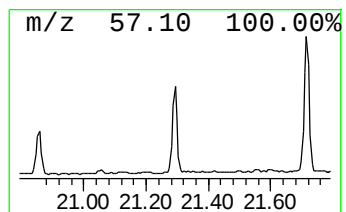
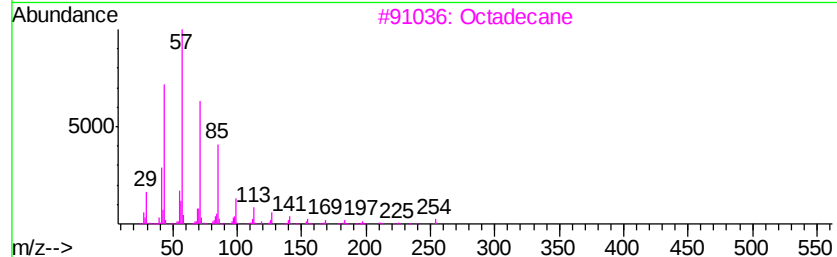
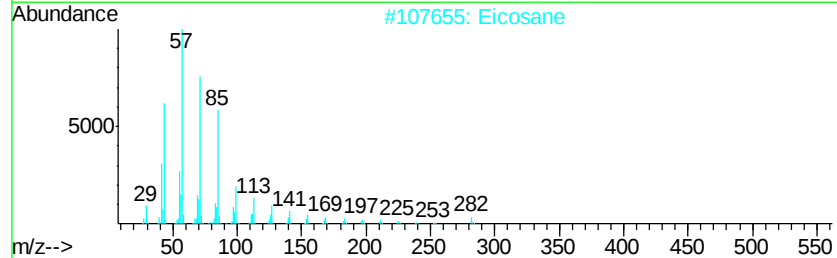
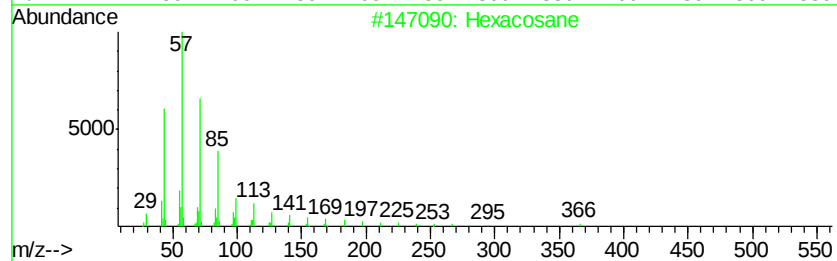
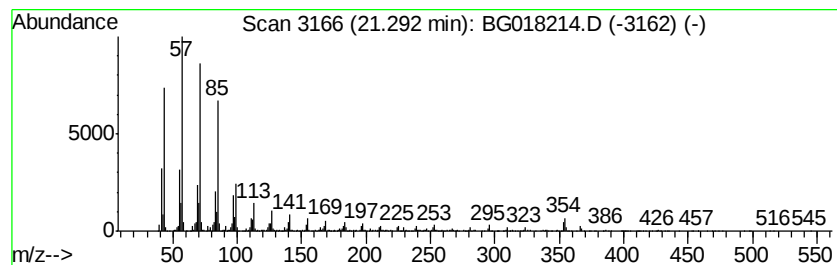
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 21.29 | 19.95 ng/ul | 880866 | Chrysene-d12     | 21.13 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexacosane              | 366 | C26H54  | 000630-01-3 | 98   |
| 2    |    | Eicosane                | 282 | C20H42  | 000112-95-8 | 97   |
| 3    |    | Octadecane              | 254 | C18H38  | 000593-45-3 | 97   |
| 4    |    | Heptadecane             | 240 | C17H36  | 000629-78-7 | 95   |
| 5    |    | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 93   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

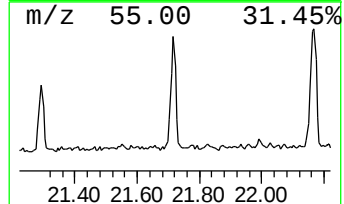
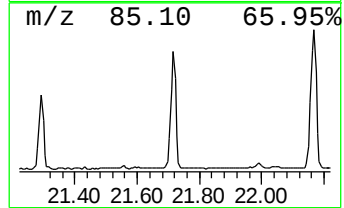
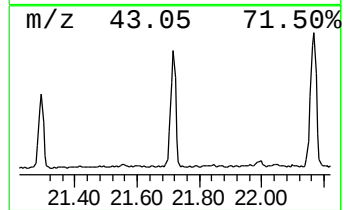
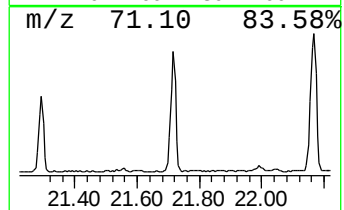
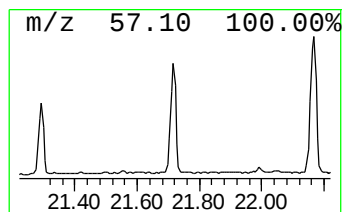
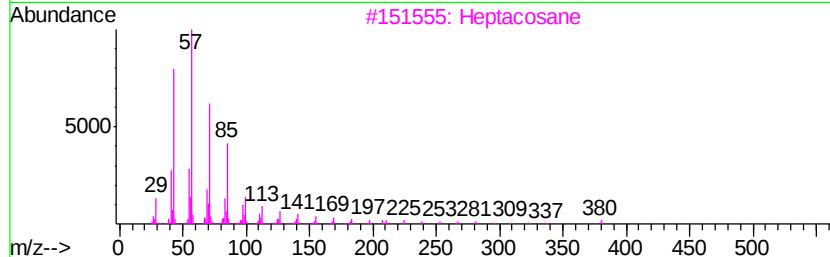
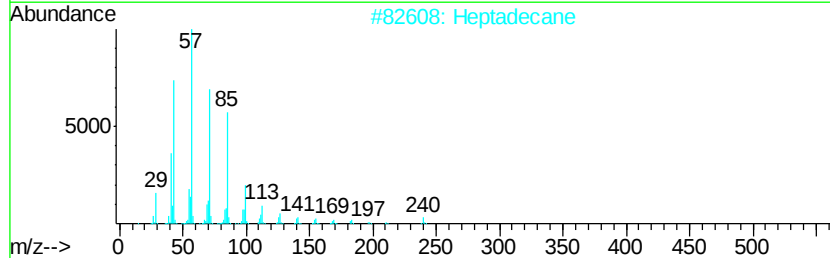
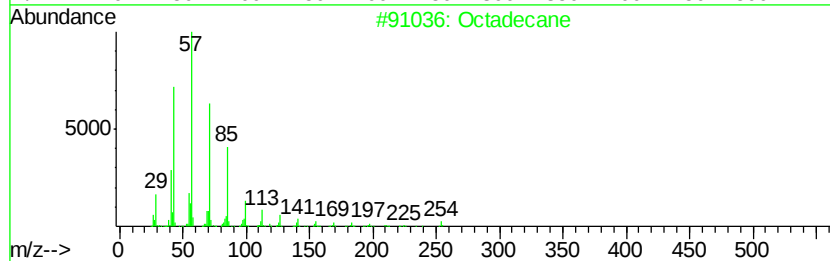
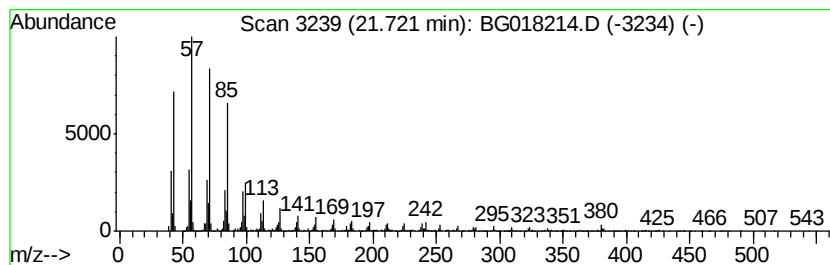
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 21.72 | 32.33 ng/ul | 1427770 | Chrysene-d12     | 21.13 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------|-----|---------|-------------|------|
| 1    | 5  | Octadecane         | 254 | C18H38  | 000593-45-3 | 96   |
| 2    |    | Heptadecane        | 240 | C17H36  | 000629-78-7 | 95   |
| 3    |    | Heptacosane        | 380 | C27H56  | 000593-49-7 | 95   |
| 4    |    | Eicosane, 9-octyl- | 394 | C28H58  | 013475-77-9 | 94   |
| 5    |    | Heptadecane        | 240 | C17H36  | 000629-78-7 | 93   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

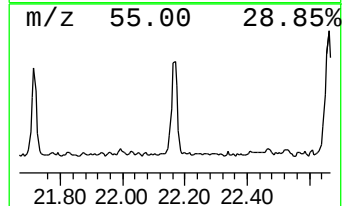
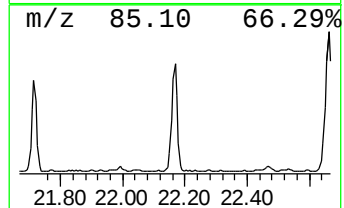
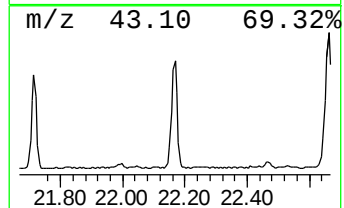
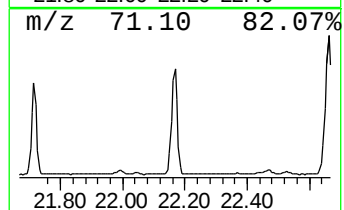
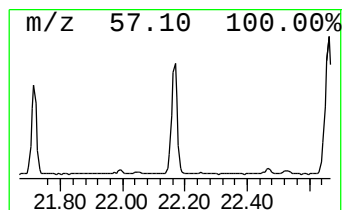
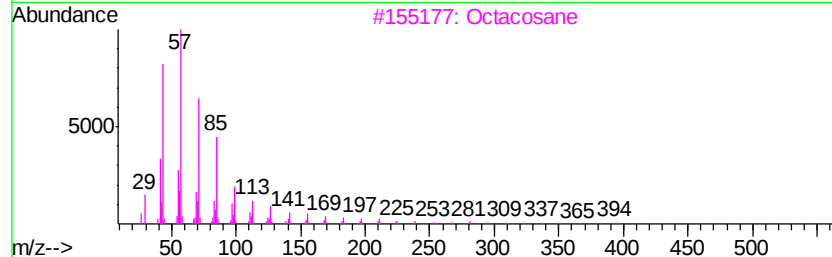
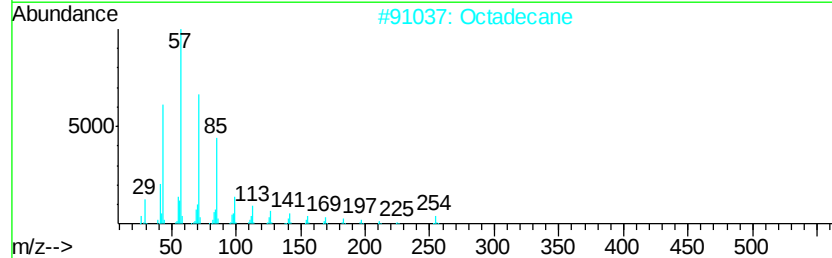
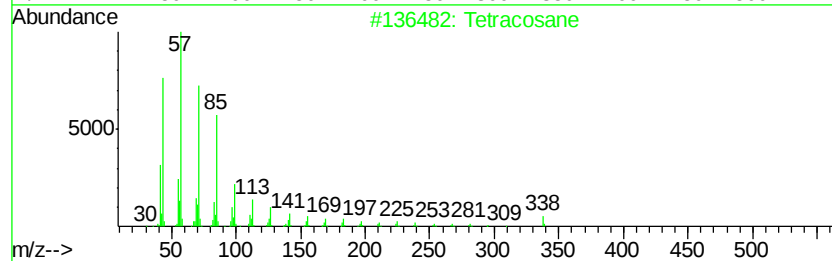
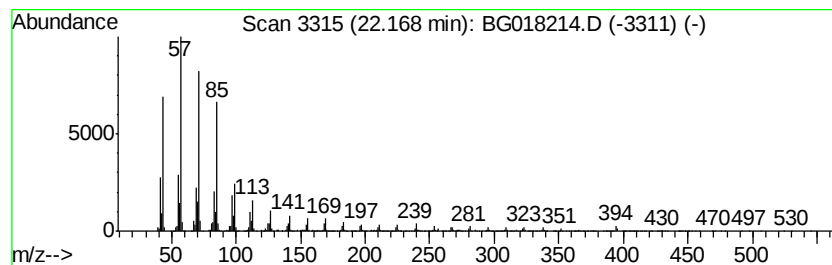
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 22.17 | 42.43 ng/ul | 1873600 | Chrysene-d12     | 21.13 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Tetracosane            | 338 | C24H50  | 000646-31-1 | 97   |
| 2    |    | Octadecane             | 254 | C18H38  | 000593-45-3 | 95   |
| 3    |    | Octacosane             | 394 | C28H58  | 000630-02-4 | 91   |
| 4    |    | Tetratetracontane      | 619 | C44H90  | 007098-22-8 | 91   |
| 5    |    | Tetracosane, 11-decyl- | 479 | C34H70  | 055429-84-0 | 91   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

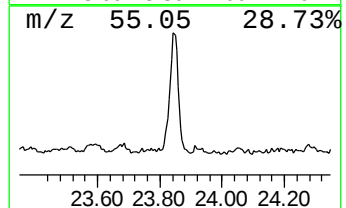
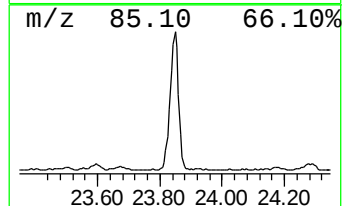
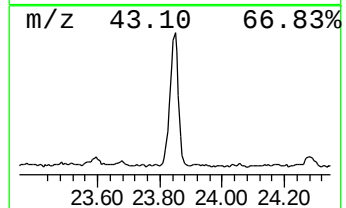
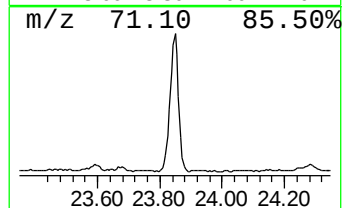
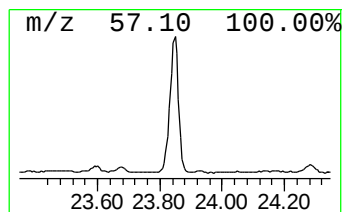
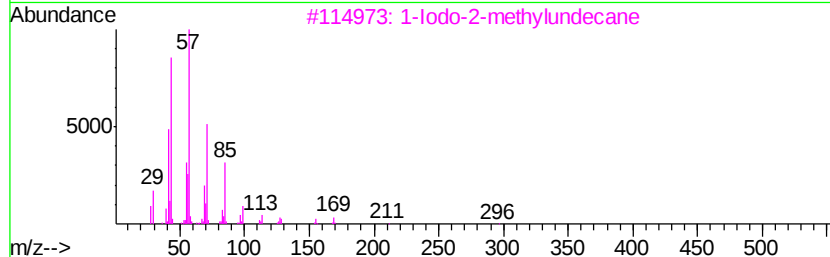
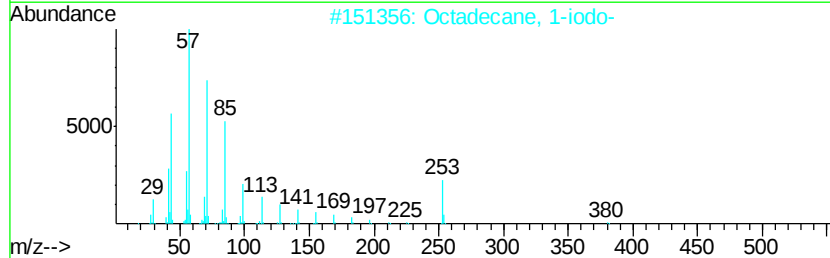
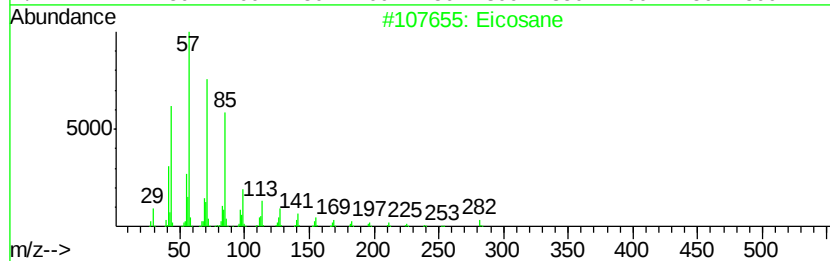
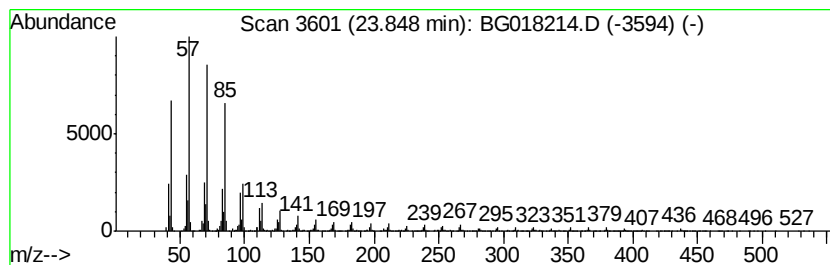
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc      | Area    | Relative to ISTD | R.T.  |
|-------|--------------|---------|------------------|-------|
| 23.85 | 123.40 ng/ul | 2680010 | Perylene-d12     | 23.31 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane                | 282 | C20H42  | 000112-95-8 | 97   |
| 2    |    | Octadecane, 1-iodo-     | 380 | C18H37I | 000629-93-6 | 94   |
| 3    |    | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 93   |
| 4    |    | Docosane, 11-decyl-     | 451 | C32H66  | 055401-55-3 | 92   |
| 5    |    | Hentriacontane          | 437 | C31H64  | 000630-04-6 | 91   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

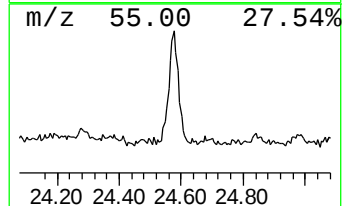
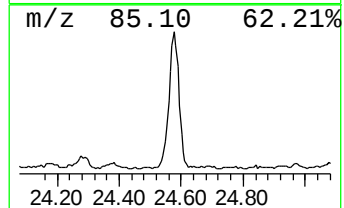
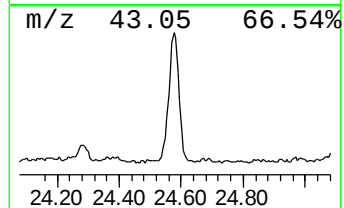
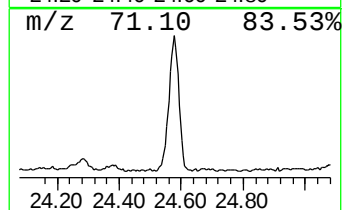
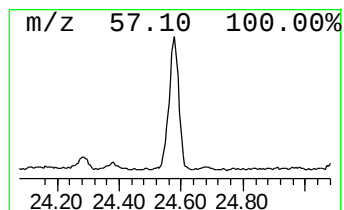
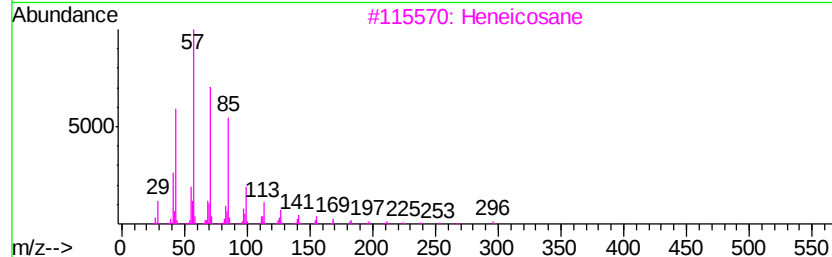
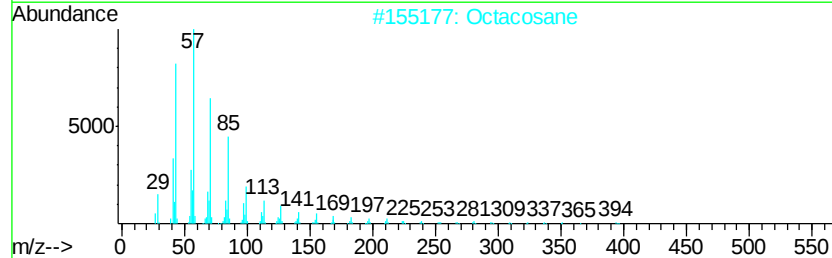
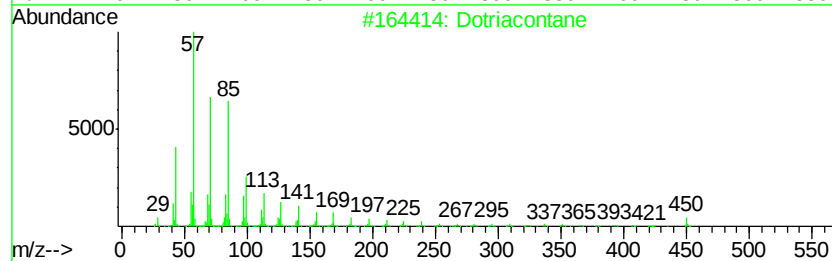
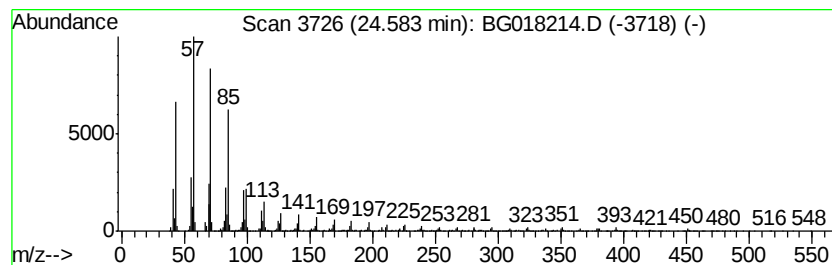
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 57 (DEL) Alkane: Cyclic24.58 Concentration Rank 2

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 24.58 | 84.18 ng/ul | 1828320 | Perylene-d12     | 23.31 |

| Hit# | of | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------|-----|----------|-------------|------|
| 1    | 5  | Dotriacontane          | 451 | C32H66   | 000544-85-4 | 97   |
| 2    |    | Octacosane             | 394 | C28H58   | 000630-02-4 | 96   |
| 3    |    | Heneicosane            | 296 | C21H44   | 000629-94-7 | 95   |
| 4    |    | Eicosane               | 282 | C20H42   | 000112-95-8 | 95   |
| 5    |    | Heptacosane, 1-chloro- | 414 | C27H55Cl | 062016-79-9 | 94   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
Data File : BG018214.D  
Acq On : 1 Aug 2015 11:03  
Operator : TP/UM  
Sample : G3073-09  
Misc :  
ALS Vial : 34 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.M  
Quant Title : SVOA CALIBRATION

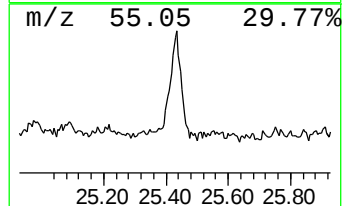
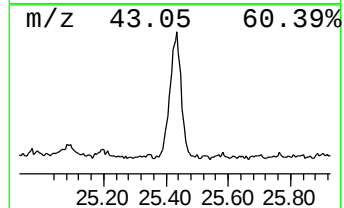
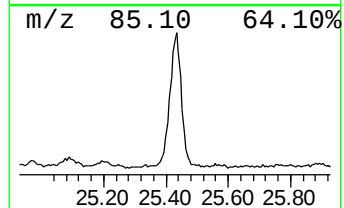
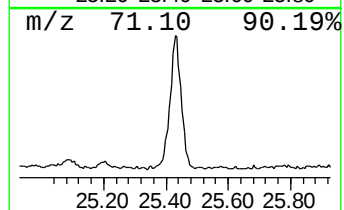
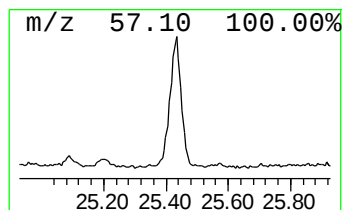
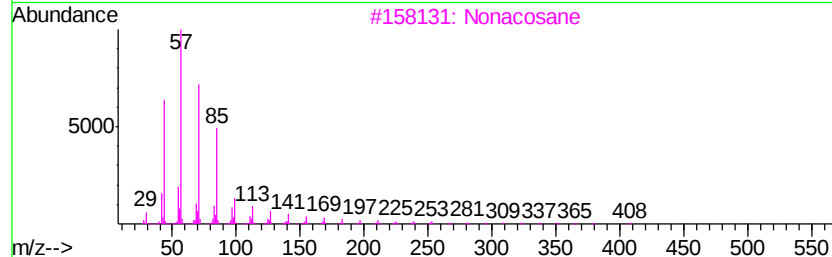
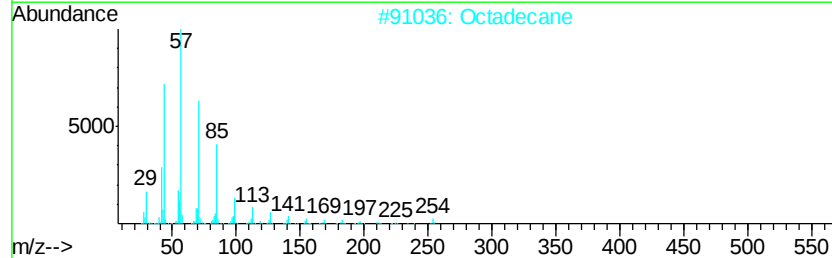
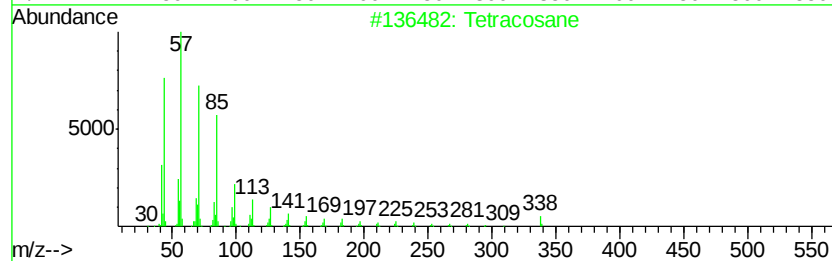
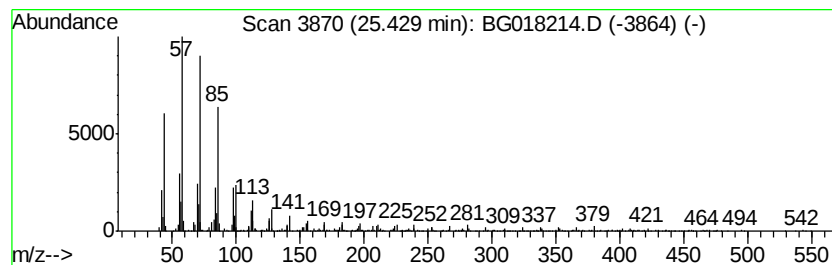
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 25.43 | 65.72 ng/ul | 1427340 | Perylene-d12     | 23.31 |

| Hit# | of | Tentative ID   | MW  | MolForm | CAS#        | Qual |
|------|----|----------------|-----|---------|-------------|------|
| 1    | 5  | Tetracosane    | 338 | C24H50  | 000646-31-1 | 97   |
| 2    |    | Octadecane     | 254 | C18H38  | 000593-45-3 | 96   |
| 3    |    | Nonacosane     | 408 | C29H60  | 000630-03-5 | 95   |
| 4    |    | Hentriacontane | 437 | C31H64  | 000630-04-6 | 93   |
| 5    |    | Tetracosane    | 338 | C24H50  | 000646-31-1 | 93   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG073115\  
 Data File : BG018214.D  
 Acq On : 1 Aug 2015 11:03  
 Operator : TP/UM  
 Sample : G3073-09  
 Misc :  
 ALS Vial : 34 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C01Q2

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG073115.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 |
| Methyl Isobutyl K... | 3.31  | 2.7     | ng/ul | 53369    | 1                     | 7.48  | 398553 | 20.0 |
| Benzene, 1-ethyl-... | 7.14  | 2.2     | ng/ul | 44116    | 1                     | 7.48  | 398553 | 20.0 |
| unknown-01           | 7.40  | 2.5     | ng/ul | 49694    | 1                     | 7.48  | 398553 | 20.0 |
| unknown-02           | 7.75  | 3.4     | ng/ul | 66711    | 1                     | 7.48  | 398553 | 20.0 |
| (DEL) Alkane: Cyc... | 8.03  | 3.3     | ng/ul | 66357    | 1                     | 7.48  | 398553 | 20.0 |
| Benzene, 1,2-diet... | 8.13  | 3.0     | ng/ul | 59427    | 1                     | 7.48  | 398553 | 20.0 |
| Naphthalene, deca... | 8.30  | 5.7     | ng/ul | 112584   | 1                     | 7.48  | 398553 | 20.0 |
| Benzene, 1-ethyl-... | 8.58  | 9.7     | ng/ul | 192711   | 1                     | 7.48  | 398553 | 20.0 |
| (DEL) Alkane: Str... | 8.80  | 4.3     | ng/ul | 86032    | 1                     | 7.48  | 398553 | 20.0 |
| unknown-03           | 8.83  | 8.0     | ng/ul | 158518   | 1                     | 7.48  | 398553 | 20.0 |
| unknown-04           | 8.94  | 4.4     | ng/ul | 170899   | 2                     | 10.27 | 769289 | 20.0 |
| (DEL) Alkane: Str... | 9.12  | 4.0     | ng/ul | 155388   | 2                     | 10.27 | 769289 | 20.0 |
| Naphthalene, deca... | 9.19  | 5.1     | ng/ul | 196404   | 2                     | 10.27 | 769289 | 20.0 |
| trans-Decalin, 2-... | 9.45  | 6.0     | ng/ul | 230631   | 2                     | 10.27 | 769289 | 20.0 |
| Benzene, 1-methyl... | 9.97  | 2.2     | ng/ul | 85661    | 2                     | 10.27 | 769289 | 20.0 |
| 2-Methyladamantane   | 10.13 | 2.5     | ng/ul | 94770    | 2                     | 10.27 | 769289 | 20.0 |
| unknown-05           | 10.68 | 3.0     | ng/ul | 114975   | 2                     | 10.27 | 769289 | 20.0 |
| 6-Tridecene, (Z)-    | 10.76 | 2.5     | ng/ul | 98004    | 2                     | 10.27 | 769289 | 20.0 |
| unknown-06           | 10.96 | 8.8     | ng/ul | 338196   | 2                     | 10.27 | 769289 | 20.0 |
| (DEL) Alkane: Str... | 11.31 | 13.9    | ng/ul | 536213   | 2                     | 10.27 | 769289 | 20.0 |
| (DEL) Alkane: Str... | 11.56 | 3.1     | ng/ul | 120979   | 2                     | 10.27 | 769289 | 20.0 |
| unknown-07           | 12.02 | 4.8     | ng/ul | 182775   | 2                     | 10.27 | 769289 | 20.0 |
| (DEL) Alkane: Cyc... | 12.43 | 5.7     | ng/ul | 199506   | 3                     | 14.17 | 698374 | 20.0 |
| unknown-08           | 12.50 | 3.8     | ng/ul | 133177   | 3                     | 14.17 | 698374 | 20.0 |
| (DEL) Alkane: Str... | 12.65 | 4.1     | ng/ul | 143209   | 3                     | 14.17 | 698374 | 20.0 |
| (DEL) Alkane: Str... | 12.70 | 15.2    | ng/ul | 530283   | 3                     | 14.17 | 698374 | 20.0 |
| unknown-09           | 12.74 | 4.6     | ng/ul | 160195   | 3                     | 14.17 | 698374 | 20.0 |
| Decahydro-4,4,8,9... | 12.93 | 6.2     | ng/ul | 214702   | 3                     | 14.17 | 698374 | 20.0 |
| (DEL) Alkane: Str... | 13.08 | 3.6     | ng/ul | 124732   | 3                     | 14.17 | 698374 | 20.0 |
| (DEL) Alkane: Str... | 14.08 | 7.3     | ng/ul | 253259   | 3                     | 14.17 | 698374 | 20.0 |
| unknown-11           | 14.71 | 9.3     | ng/ul | 324420   | 3                     | 14.17 | 698374 | 20.0 |
| unknown-12           | 14.95 | 8.1     | ng/ul | 282509   | 3                     | 14.17 | 698374 | 20.0 |
| (DEL) Alkane: Str... | 15.04 | 4.8     | ng/ul | 168341   | 3                     | 14.17 | 698374 | 20.0 |
| (DEL) Alkane: Str... | 15.45 | 13.3    | ng/ul | 464960   | 3                     | 14.17 | 698374 | 20.0 |
| Cyclohexane, 1,1'... | 15.65 | 6.1     | ng/ul | 154240   | 4                     | 16.92 | 506624 | 20.0 |
| (DEL) Alkane: Str... | 15.93 | 42.8    | ng/ul | 1083290  | 4                     | 16.92 | 506624 | 20.0 |
| (DEL) Alkane: Str... | 16.70 | 4.3     | ng/ul | 107870   | 4                     | 16.92 | 506624 | 20.0 |
| (DEL) Alkane: Str... | 16.74 | 19.3    | ng/ul | 488722   | 4                     | 16.92 | 506624 | 20.0 |
| (DEL) Alkane: Str... | 17.43 | 4.6     | ng/ul | 116102   | 4                     | 16.92 | 506624 | 20.0 |
| unknown-13           | 17.84 | 5.0     | ng/ul | 126470   | 4                     | 16.92 | 506624 | 20.0 |
| 4b,8-Dimethyl-2-i... | 18.55 | 12.3    | ng/ul | 310829   | 4                     | 16.92 | 506624 | 20.0 |
| 10,18-Bisnorabiet... | 19.05 | 8.9     | ng/ul | 394643   | 5                     | 21.13 | 883175 | 20.0 |
| (DEL) Alkane: Str... | 20.39 | 3.4     | ng/ul | 150760   | 5                     | 21.13 | 883175 | 20.0 |
| (DEL) Alkane: Str... | 20.86 | 10.9    | ng/ul | 478938   | 5                     | 21.13 | 883175 | 20.0 |
| (DEL) Alkane: Str... | 21.29 | 19.9    | ng/ul | 880866   | 5                     | 21.13 | 883175 | 20.0 |
| (DEL) Alkane: Str... | 21.72 | 32.3    | ng/ul | 1427770  | 5                     | 21.13 | 883175 | 20.0 |
| (DEL) Alkane: Str... | 22.17 | 42.4    | ng/ul | 1873600  | 5                     | 21.13 | 883175 | 20.0 |
| (DEL) Alkane: Str... | 23.85 | 123.4   | ng/ul | 2680010  | 6                     | 23.31 | 434379 | 20.0 |
| (DEL) Alkane: Cyc... | 24.58 | 84.2    | ng/ul | 1828320  | 6                     | 23.31 | 434379 | 20.0 |

(DEL) Alkane: Str... 25.43 65.7 ng/ul 1427340 6 23.31 434379 20.0

Instrument :

BNA\_G

ClientSampleId :

C01Q2