

Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
 Data File : BG018970.D  
 Acq On : 30 Sep 2015 13:28  
 Operator : UM/NP  
 Sample : G3690-01DL 5X  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BC2W1DL

## 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-BG091015.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         | 3.156       | 49            | 54          | 69           | rVB      | 431879         | 590974        | 1.25%           | 0.475%        |
| 2         | 5.624       | 468           | 474         | 484          | rBV      | 231705         | 459035        | 0.97%           | 0.369%        |
| 3         | 5.994       | 531           | 537         | 545          | rVB      | 207878         | 333257        | 0.71%           | 0.268%        |
| 4         | 6.975       | 699           | 704         | 709          | rVB      | 164590         | 243552        | 0.52%           | 0.196%        |
| 5         | 7.134       | 725           | 731         | 737          | rBV5     | 145216         | 327809        | 0.69%           | 0.263%        |
| 6         | 7.204       | 737           | 743         | 750          | rVB3     | 90202          | 196709        | 0.42%           | 0.158%        |
| 7         | 7.604       | 805           | 811         | 815          | rBV      | 394858         | 627929        | 1.33%           | 0.504%        |
| 8         | 7.909       | 853           | 863         | 866          | rBV5     | 51129          | 160167        | 0.34%           | 0.129%        |
| 9         | 7.962       | 866           | 872         | 881          | rVV3     | 326950         | 758717        | 1.61%           | 0.609%        |
| 10        | 8.109       | 891           | 897         | 903          | rBV3     | 156108         | 291790        | 0.62%           | 0.234%        |
| 11        | 8.215       | 910           | 915         | 921          | rVV      | 395568         | 730274        | 1.55%           | 0.587%        |
| 12        | 8.267       | 921           | 924         | 932          | rVB3     | 98396          | 173809        | 0.37%           | 0.140%        |
| 13        | 8.526       | 962           | 968         | 973          | rBV2     | 352325         | 646849        | 1.37%           | 0.520%        |
| 14        | 8.649       | 980           | 989         | 995          | rBV3     | 234054         | 621934        | 1.32%           | 0.500%        |
| 15        | 8.861       | 1021          | 1025        | 1029         | rBV      | 105981         | 166921        | 0.35%           | 0.134%        |
| 16        | 9.090       | 1059          | 1064        | 1070         | rVB3     | 143597         | 257686        | 0.55%           | 0.207%        |
| 17        | 9.208       | 1070          | 1084        | 1095         | rVB      | 447411         | 1001712       | 2.12%           | 0.805%        |
| 18        | 9.337       | 1102          | 1106        | 1115         | rVB7     | 80441          | 192788        | 0.41%           | 0.155%        |
| 19        | 9.971       | 1208          | 1214        | 1222         | rBV5     | 63573          | 147348        | 0.31%           | 0.118%        |
| 20        | 10.189      | 1246          | 1251        | 1258         | rVV5     | 101811         | 192316        | 0.41%           | 0.154%        |
| 21        | 10.659      | 1324          | 1331        | 1342         | rBV      | 476585         | 952422        | 2.02%           | 0.765%        |
| 22        | 10.753      | 1342          | 1347        | 1349         | rBV      | 139000         | 211845        | 0.45%           | 0.170%        |
| 23        | 10.788      | 1349          | 1353        | 1360         | rVB      | 314024         | 541673        | 1.15%           | 0.435%        |
| 24        | 12.122      | 1575          | 1580        | 1587         | rVB      | 126069         | 199577        | 0.42%           | 0.160%        |
| 25        | 12.192      | 1587          | 1592        | 1601         | rVV      | 111037         | 169942        | 0.36%           | 0.137%        |
| 26        | 13.127      | 1746          | 1751        | 1760         | rVB4     | 157928         | 327540        | 0.69%           | 0.263%        |
| 27        | 13.432      | 1797          | 1803        | 1815         | rVV2     | 120734         | 243558        | 0.52%           | 0.196%        |
| 28        | 13.638      | 1833          | 1838        | 1840         | rVV      | 115154         | 174155        | 0.37%           | 0.140%        |
| 29        | 13.661      | 1840          | 1842        | 1852         | rVB2     | 116869         | 178469        | 0.38%           | 0.143%        |
| 30        | 14.255      | 1938          | 1943        | 1946         | rBV      | 102185         | 161716        | 0.34%           | 0.130%        |
| 31        | 14.307      | 1949          | 1952        | 1956         | rVV      | 88442          | 138828        | 0.29%           | 0.112%        |
| 32        | 14.372      | 1956          | 1963        | 1971         | rVV4     | 84527          | 207050        | 0.44%           | 0.166%        |
| 33        | 14.501      | 1982          | 1985        | 1993         | rVB      | 138606         | 176861        | 0.37%           | 0.142%        |
| 34        | 14.613      | 1999          | 2004        | 2010         | rVB2     | 625509         | 1029311       | 2.18%           | 0.827%        |

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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-BG091015.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |          |          |         |         |
|----|--------|------|------|------|------|----------|----------|---------|---------|
| 35 | 14.695 | 2010 | 2018 | 2023 | rVB2 | 70161    | 139796   | 0.30%   | 0.112%  |
| 36 | 14.895 | 2044 | 2052 | 2062 | rVV  | 3936912  | 6388835  | 13.52%  | 5.132%  |
| 37 | 15.036 | 2072 | 2076 | 2086 | rVB2 | 109642   | 177354   | 0.38%   | 0.142%  |
| 38 | 15.206 | 2101 | 2105 | 2107 | rVV  | 215953   | 316270   | 0.67%   | 0.254%  |
| 39 | 15.242 | 2107 | 2111 | 2117 | rVV  | 1393526  | 2129467  | 4.51%   | 1.711%  |
| 40 | 15.447 | 2142 | 2146 | 2152 | rVB  | 98134    | 147098   | 0.31%   | 0.118%  |
| 41 | 15.612 | 2168 | 2174 | 2179 | rVB  | 957635   | 1412839  | 2.99%   | 1.135%  |
| 42 | 15.747 | 2191 | 2197 | 2207 | rBV  | 862105   | 1389253  | 2.94%   | 1.116%  |
| 43 | 16.146 | 2261 | 2265 | 2269 | rBV2 | 274586   | 326583   | 0.69%   | 0.262%  |
| 44 | 16.311 | 2288 | 2293 | 2296 | rBV  | 278623   | 413367   | 0.87%   | 0.332%  |
| 45 | 16.340 | 2296 | 2298 | 2309 | rVB6 | 204193   | 368702   | 0.78%   | 0.296%  |
| 46 | 16.434 | 2310 | 2314 | 2318 | rBV  | 137571   | 196154   | 0.42%   | 0.158%  |
| 47 | 16.493 | 2320 | 2324 | 2329 | rBV2 | 140476   | 216255   | 0.46%   | 0.174%  |
| 48 | 16.552 | 2329 | 2334 | 2338 | rBV3 | 200690   | 288350   | 0.61%   | 0.232%  |
| 49 | 16.763 | 2365 | 2370 | 2378 | rBV5 | 177648   | 340418   | 0.72%   | 0.273%  |
| 50 | 17.069 | 2407 | 2422 | 2426 | rBV  | 16047649 | 47249873 | 100.00% | 37.955% |
| 51 | 17.157 | 2434 | 2437 | 2446 | rVB2 | 146920   | 265043   | 0.56%   | 0.213%  |
| 52 | 17.369 | 2468 | 2473 | 2478 | rVB2 | 806223   | 1185265  | 2.51%   | 0.952%  |
| 53 | 17.433 | 2480 | 2484 | 2487 | rVV3 | 133687   | 221715   | 0.47%   | 0.178%  |
| 54 | 17.463 | 2487 | 2489 | 2495 | rVB2 | 151932   | 205039   | 0.43%   | 0.165%  |
| 55 | 17.598 | 2509 | 2512 | 2517 | rVB2 | 102911   | 134570   | 0.28%   | 0.108%  |
| 56 | 17.709 | 2526 | 2531 | 2537 | rBV  | 1012455  | 1553989  | 3.29%   | 1.248%  |
| 57 | 17.762 | 2538 | 2540 | 2544 | rVB2 | 135958   | 181591   | 0.38%   | 0.146%  |
| 58 | 17.845 | 2550 | 2554 | 2564 | rVB2 | 384006   | 718348   | 1.52%   | 0.577%  |
| 59 | 18.062 | 2587 | 2591 | 2596 | rVB3 | 254098   | 359732   | 0.76%   | 0.289%  |
| 60 | 18.273 | 2621 | 2627 | 2630 | rBV2 | 1220622  | 1779678  | 3.77%   | 1.430%  |
| 61 | 18.320 | 2630 | 2635 | 2640 | rVB  | 2983937  | 4189007  | 8.87%   | 3.365%  |
| 62 | 18.367 | 2640 | 2643 | 2648 | rBV2 | 122841   | 241050   | 0.51%   | 0.194%  |
| 63 | 18.438 | 2652 | 2655 | 2660 | rVB3 | 368200   | 621461   | 1.32%   | 0.499%  |
| 64 | 18.550 | 2668 | 2674 | 2678 | rBV2 | 607767   | 1058173  | 2.24%   | 0.850%  |
| 65 | 18.661 | 2690 | 2693 | 2694 | rBV3 | 118806   | 136748   | 0.29%   | 0.110%  |
| 66 | 18.691 | 2694 | 2698 | 2703 | rVB3 | 514453   | 755061   | 1.60%   | 0.607%  |
| 67 | 18.785 | 2711 | 2714 | 2718 | rBV2 | 137428   | 181414   | 0.38%   | 0.146%  |
| 68 | 18.931 | 2736 | 2739 | 2742 | rBV4 | 149456   | 212388   | 0.45%   | 0.171%  |
| 69 | 18.967 | 2742 | 2745 | 2748 | rVV2 | 137784   | 177076   | 0.37%   | 0.142%  |
| 70 | 19.090 | 2763 | 2766 | 2770 | rBV4 | 123687   | 197221   | 0.42%   | 0.158%  |
| 71 | 19.161 | 2774 | 2778 | 2782 | rVB  | 896262   | 995141   | 2.11%   | 0.799%  |

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

Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG091015.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |       |         |         |       |        |
|-----|--------|------|------|------|-------|---------|---------|-------|--------|
| 72  | 19.431 | 2817 | 2824 | 2833 | rBV2  | 1494954 | 3946756 | 8.35% | 3.170% |
| 73  | 19.543 | 2839 | 2843 | 2850 | rBV2  | 405807  | 546455  | 1.16% | 0.439% |
| 74  | 19.607 | 2851 | 2854 | 2859 | rVV   | 484687  | 587529  | 1.24% | 0.472% |
| 75  | 19.736 | 2872 | 2876 | 2881 | rVB2  | 1369935 | 1513780 | 3.20% | 1.216% |
| 76  | 19.795 | 2882 | 2886 | 2891 | rVB   | 353731  | 470840  | 1.00% | 0.378% |
| 77  | 20.118 | 2937 | 2941 | 2945 | rVB   | 801744  | 962069  | 2.04% | 0.773% |
| 78  | 20.265 | 2962 | 2966 | 2973 | rVB   | 1713246 | 2023357 | 4.28% | 1.625% |
| 79  | 20.577 | 3015 | 3019 | 3023 | rBV   | 495571  | 668296  | 1.41% | 0.537% |
| 80  | 20.624 | 3023 | 3027 | 3033 | rVV   | 554814  | 747132  | 1.58% | 0.600% |
| 81  | 20.706 | 3037 | 3041 | 3045 | rVV   | 737234  | 967087  | 2.05% | 0.777% |
| 82  | 20.759 | 3045 | 3050 | 3058 | rVB   | 1817326 | 2253105 | 4.77% | 1.810% |
| 83  | 20.994 | 3087 | 3090 | 3094 | rBV   | 179469  | 226775  | 0.48% | 0.182% |
| 84  | 21.111 | 3106 | 3110 | 3114 | rBV   | 322486  | 463829  | 0.98% | 0.373% |
| 85  | 21.258 | 3130 | 3135 | 3143 | rVB   | 1651027 | 2246501 | 4.75% | 1.805% |
| 86  | 21.499 | 3172 | 3176 | 3181 | rBV   | 618028  | 784176  | 1.66% | 0.630% |
| 87  | 21.622 | 3192 | 3197 | 3210 | rVB   | 947624  | 1730010 | 3.66% | 1.390% |
| 88  | 21.799 | 3222 | 3227 | 3233 | rVB   | 1352908 | 1802573 | 3.81% | 1.448% |
| 89  | 22.233 | 3297 | 3301 | 3307 | rVB   | 367010  | 601902  | 1.27% | 0.484% |
| 90  | 22.398 | 3323 | 3329 | 3338 | rVB   | 1057600 | 1667809 | 3.53% | 1.340% |
| 91  | 22.756 | 3385 | 3390 | 3396 | rBV   | 224819  | 426898  | 0.90% | 0.343% |
| 92  | 23.080 | 3439 | 3445 | 3453 | rVB   | 735317  | 1359258 | 2.88% | 1.092% |
| 93  | 23.867 | 3572 | 3579 | 3588 | rBV   | 542252  | 1157103 | 2.45% | 0.929% |
| 94  | 24.801 | 3730 | 3738 | 3754 | rVB3  | 767216  | 2294258 | 4.86% | 1.843% |
| 95  | 25.077 | 3778 | 3785 | 3788 | rBV5  | 88289   | 228845  | 0.48% | 0.184% |
| 96  | 25.906 | 3919 | 3926 | 3937 | rVB4  | 215255  | 611071  | 1.29% | 0.491% |
| 97  | 26.511 | 4021 | 4029 | 4052 | rVB10 | 202910  | 978747  | 2.07% | 0.786% |
| 98  | 27.228 | 4140 | 4151 | 4164 | rVB4  | 340645  | 1212118 | 2.57% | 0.974% |
| 99  | 27.375 | 4165 | 4176 | 4192 | rVB4  | 354409  | 1291004 | 2.73% | 1.037% |
| 100 | 28.579 | 4369 | 4381 | 4400 | rVB3  | 225346  | 1014307 | 2.15% | 0.815% |

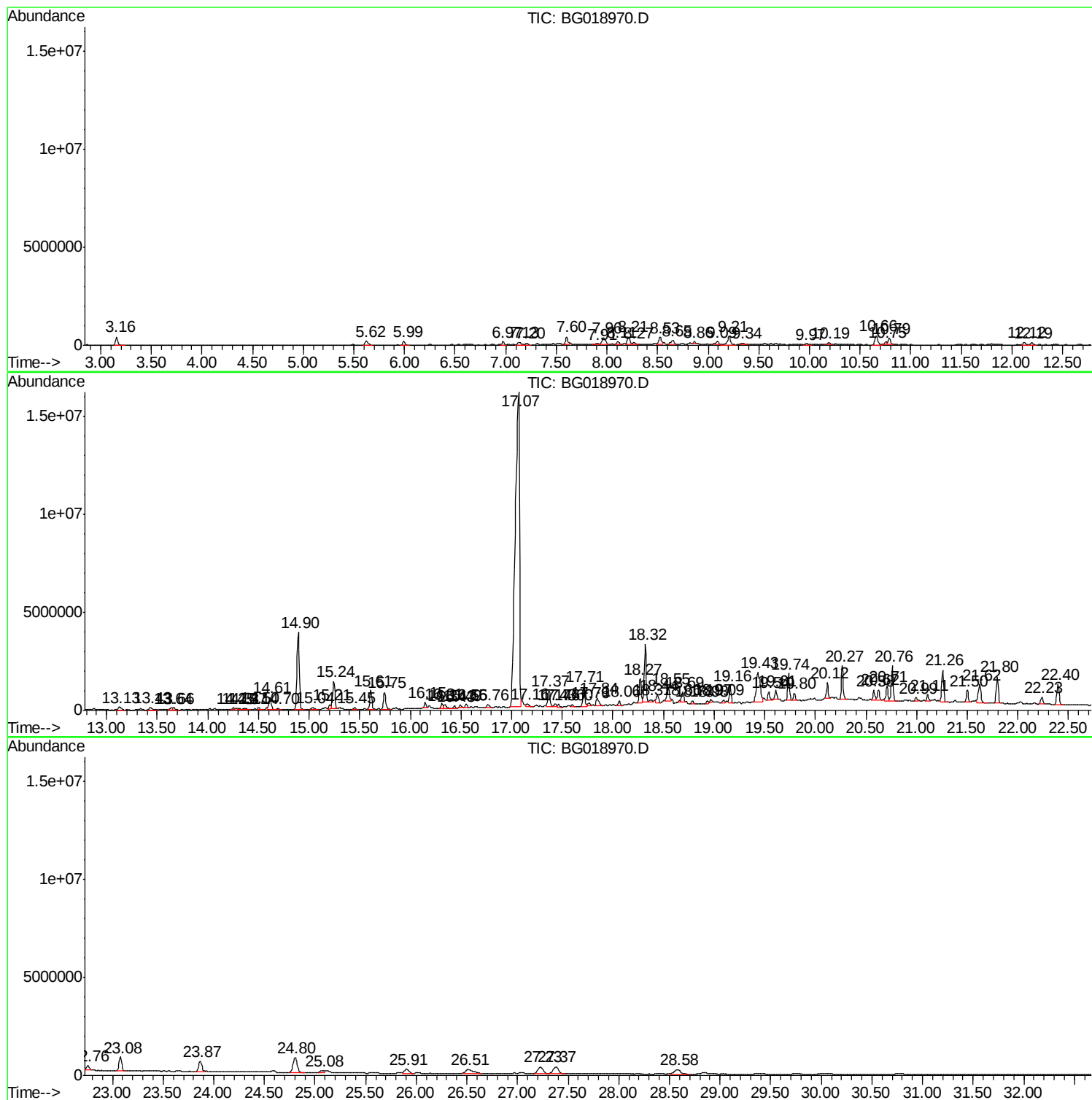
Sum of corrected areas: 124488437

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

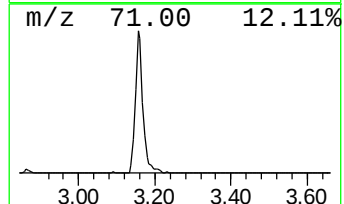
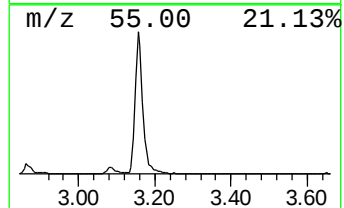
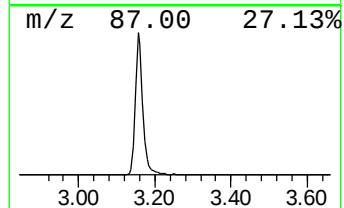
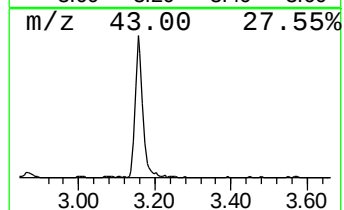
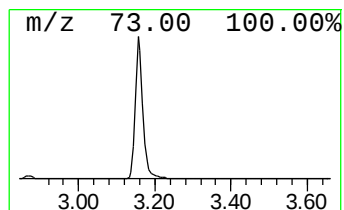
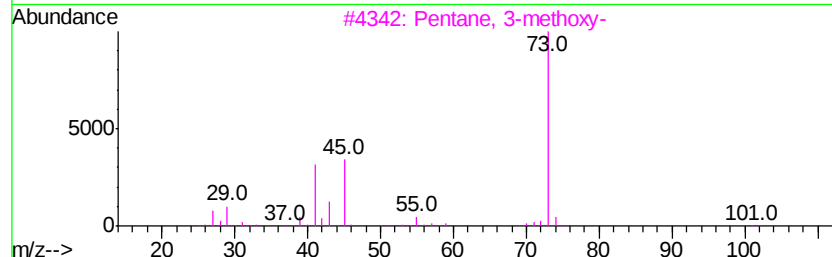
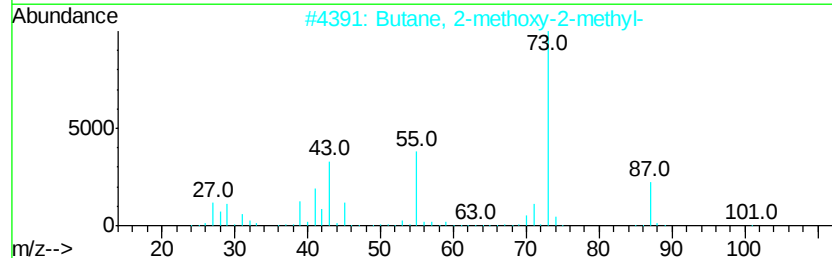
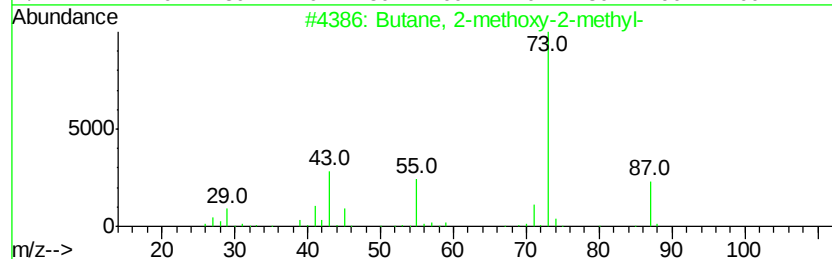
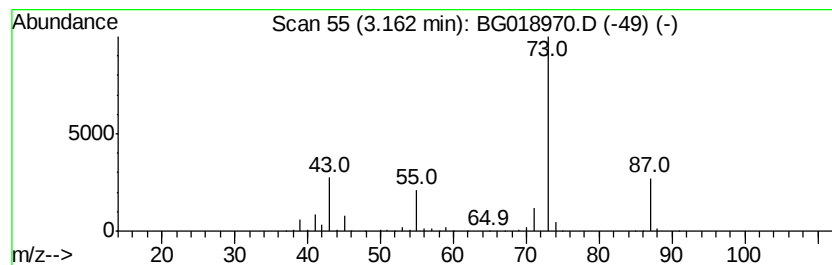
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 17

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 3.16 | 15.58 ng/ul | 590974 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 78   |
| 3    |    |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 4    |    |   | Acetamide, N-ethyl-         | 87  | C4H9NO  | 000625-50-3 | 9    |
| 5    |    |   | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |



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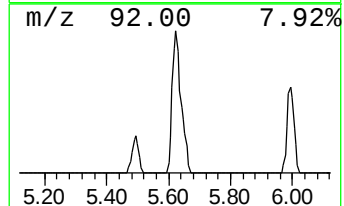
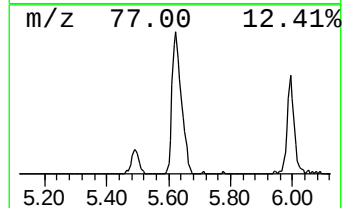
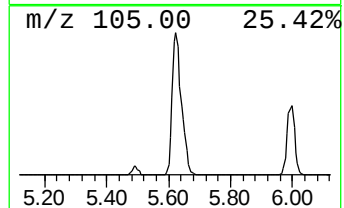
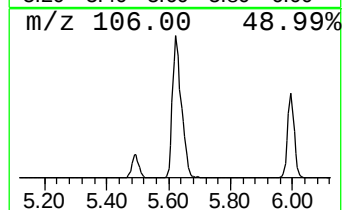
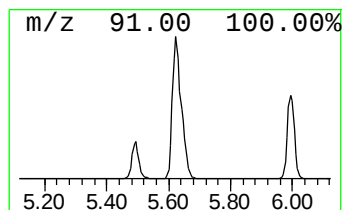
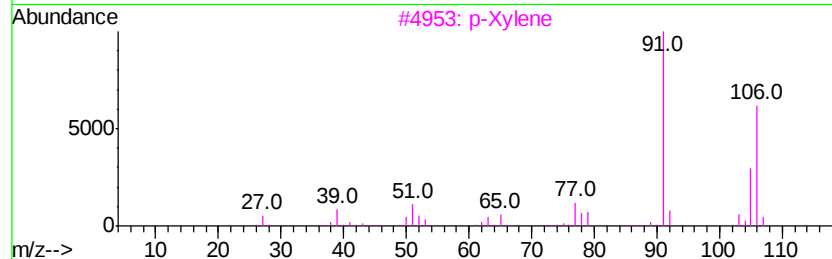
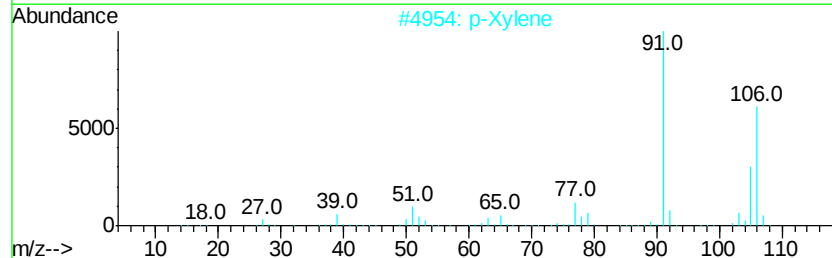
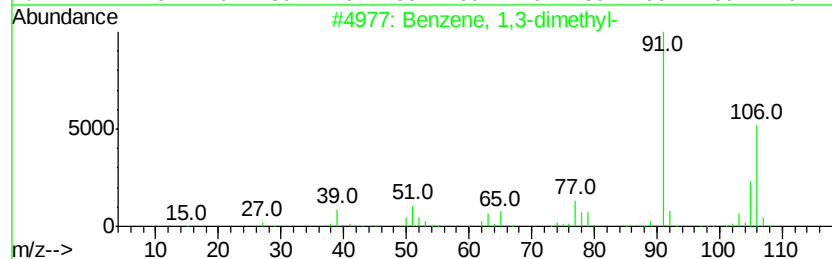
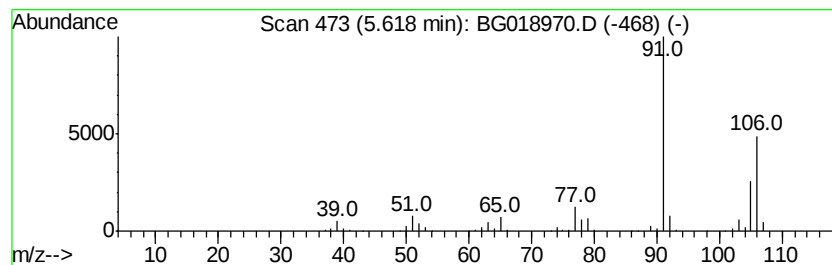
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 Benzene, 1,3-dimethyl- Concentration Rank 20

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 5.62 | 12.10 ng/ul | 459035 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 97   |
| 2    |    | p-Xylene               | 106 | C8H10   | 000106-42-3 | 97   |
| 3    |    | p-Xylene               | 106 | C8H10   | 000106-42-3 | 97   |
| 4    |    | o-Xylene               | 106 | C8H10   | 000095-47-6 | 95   |
| 5    |    | p-Xylene               | 106 | C8H10   | 000106-42-3 | 95   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

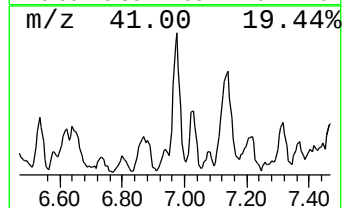
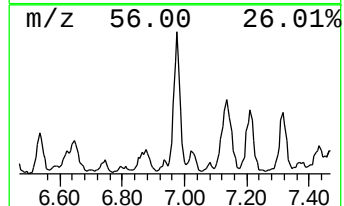
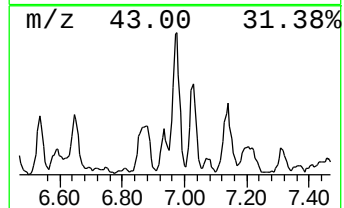
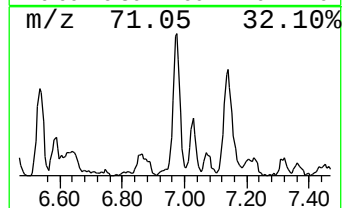
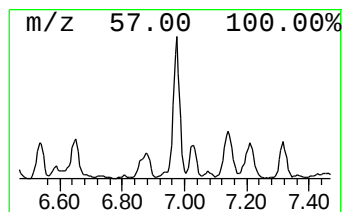
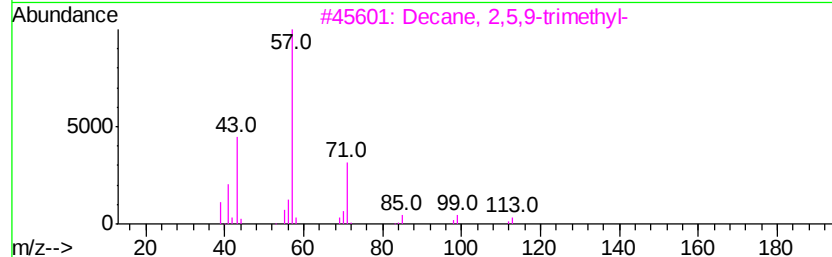
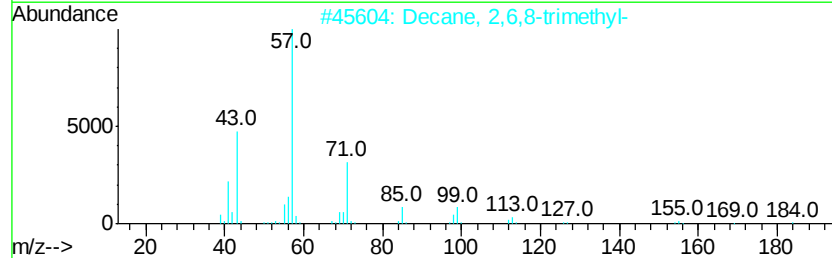
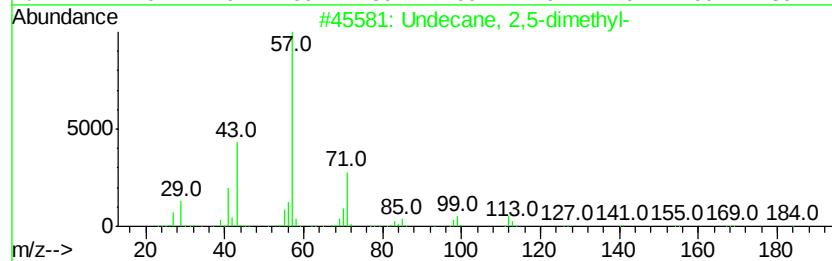
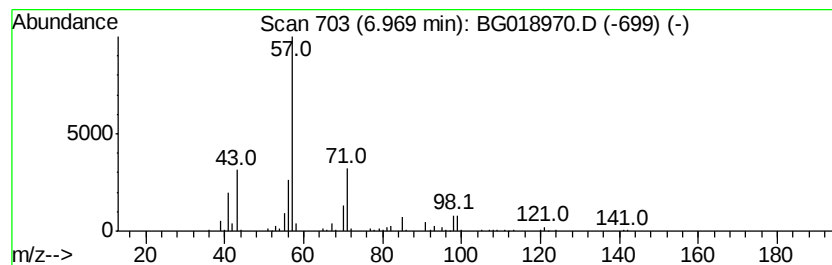
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 39

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.97 | 6.42 ng/ul | 243552 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Undecane, 2,5-dimethyl-  | 184 | C13H28  | 017301-22-3 | 59   |
| 2    |    |   | Decane, 2,6,8-trimethyl- | 184 | C13H28  | 062108-26-3 | 59   |
| 3    |    |   | Decane, 2,5,9-trimethyl- | 184 | C13H28  | 062108-22-9 | 59   |
| 4    |    |   | Heptane, 2,5-dimethyl-   | 128 | C9H20   | 002216-30-0 | 53   |
| 5    |    |   | Dodecane, 6-methyl-      | 184 | C13H28  | 006044-71-9 | 53   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

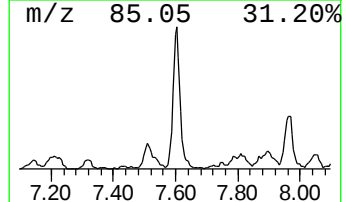
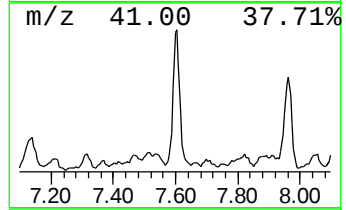
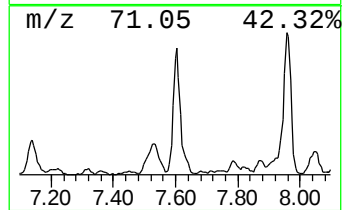
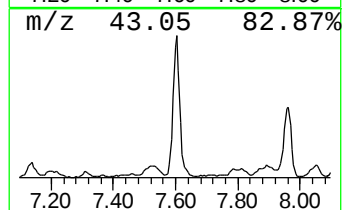
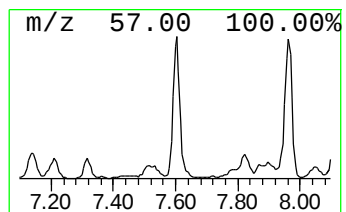
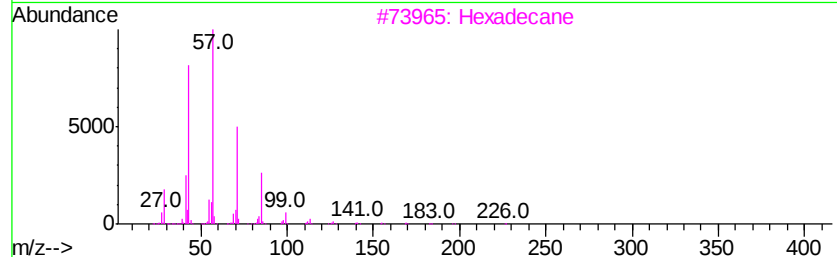
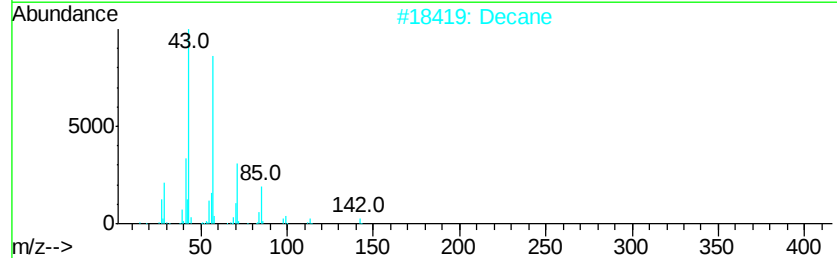
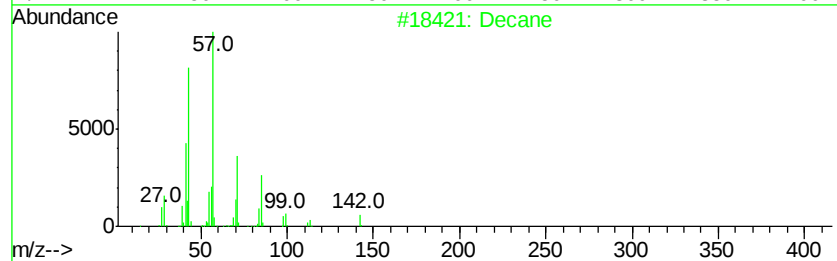
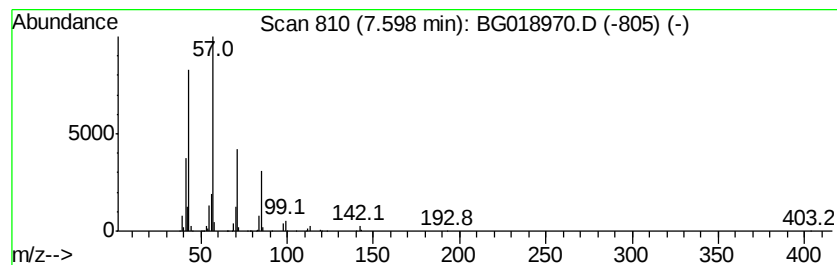
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.60 | 16.55 ng/ul | 627929 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Decane       | 142 | C10H22  | 000124-18-5 | 95   |
| 2    |    | Decane       | 142 | C10H22  | 000124-18-5 | 94   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 4    |    | Decane       | 142 | C10H22  | 000124-18-5 | 90   |
| 5    |    | Pentadecane  | 212 | C15H32  | 000629-62-9 | 78   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

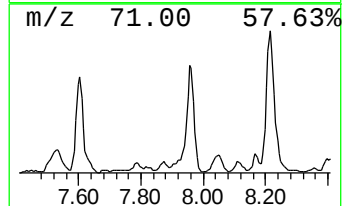
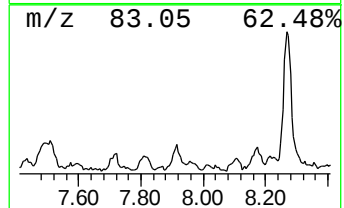
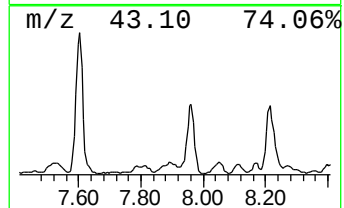
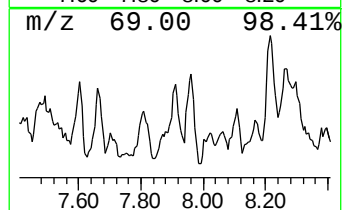
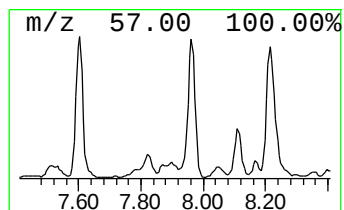
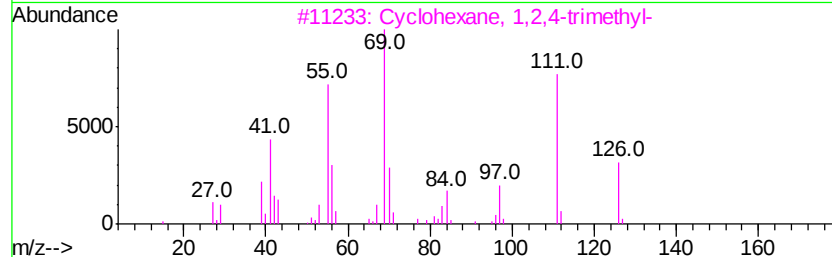
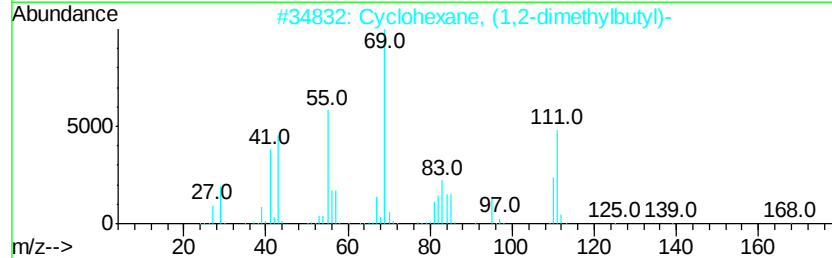
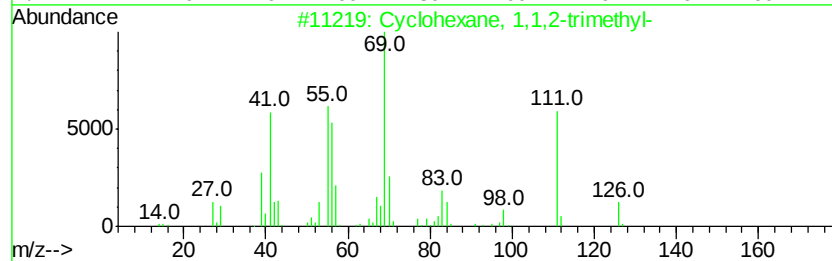
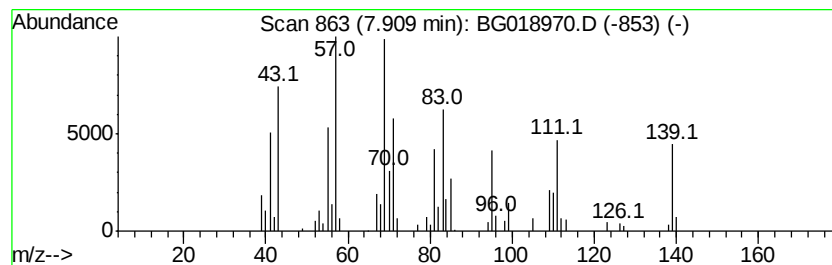
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Cyclic7.91 Concentration Rank 56

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.91 | 4.22 ng/ul | 160167 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,1,2-trimethyl-     | 126 | C9H18   | 007094-26-0 | 38   |
| 2    |    |   | Cyclohexane, (1,2-dimethylbutyl)- | 168 | C12H24  | 061142-37-8 | 35   |
| 3    |    |   | Cyclohexane, 1,2,4-trimethyl-     | 126 | C9H18   | 002234-75-5 | 30   |
| 4    |    |   | 1,1'-Bicyclooctyl                 | 222 | C16H30  | 006708-17-4 | 27   |
| 5    |    |   | 3-Octene, 2,2-dimethyl-           | 140 | C10H20  | 086869-76-3 | 25   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

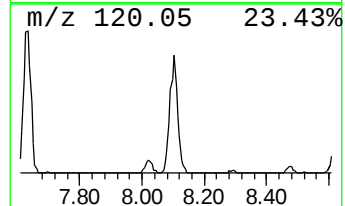
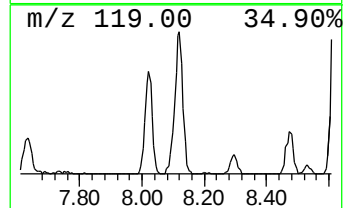
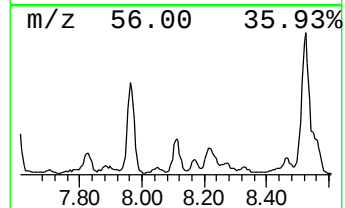
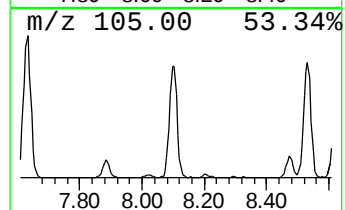
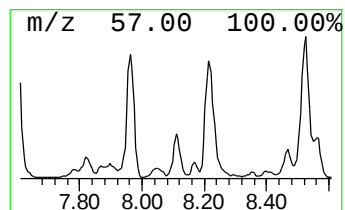
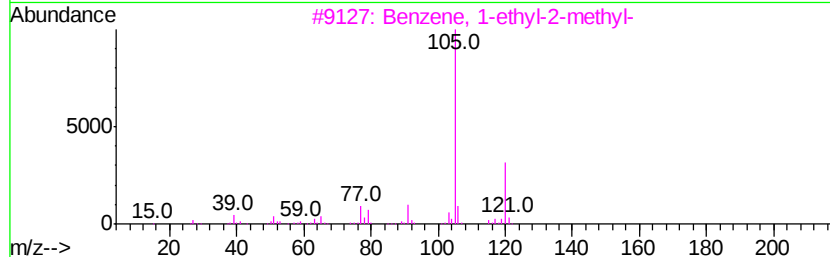
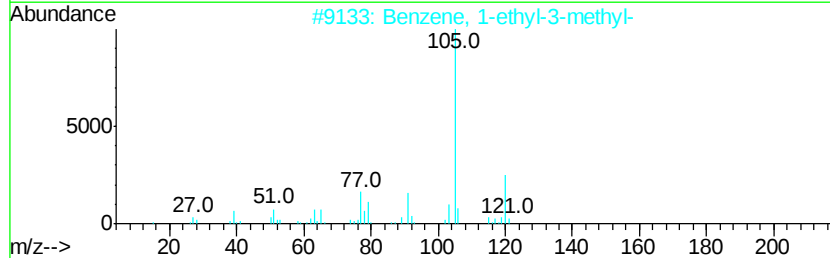
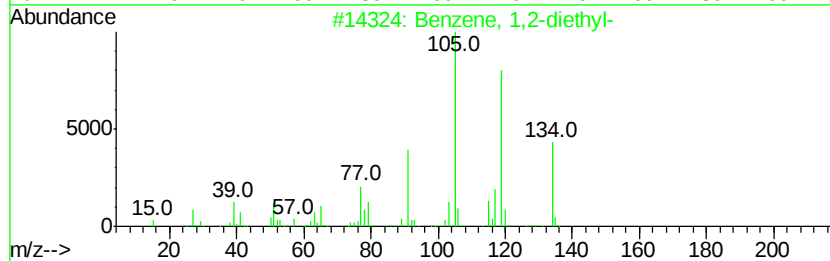
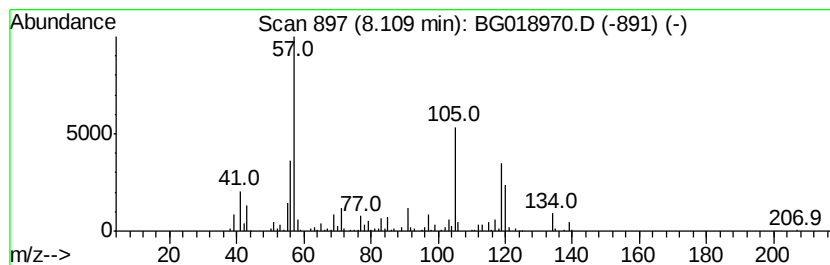
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 unknown-01 Concentration Rank 32

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.11 | 7.69 ng/ul | 291790 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2-diethyl-      | 134 | C10H14  | 000135-01-3 | 43   |
| 2    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 38   |
| 3    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 38   |
| 4    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 30   |
| 5    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 30   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

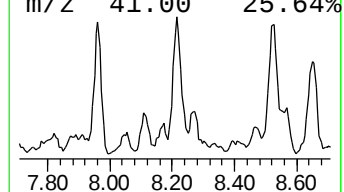
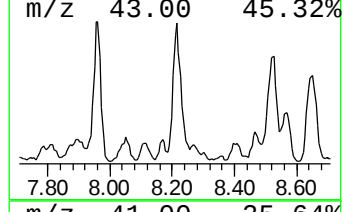
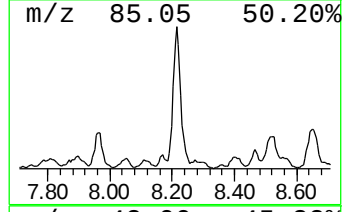
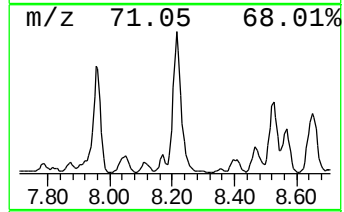
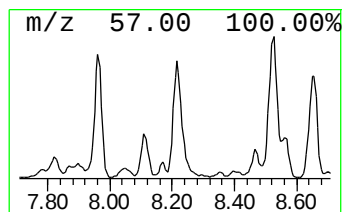
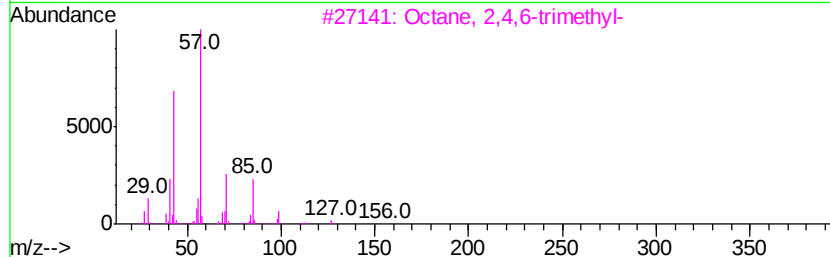
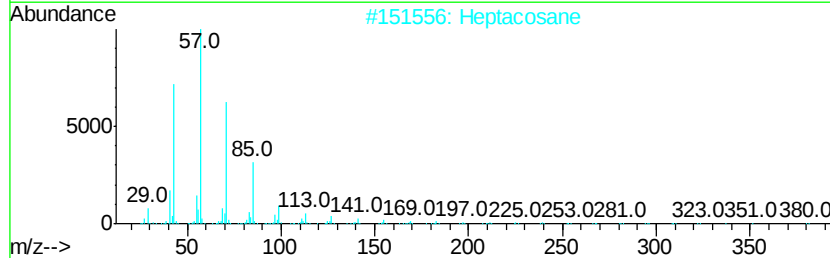
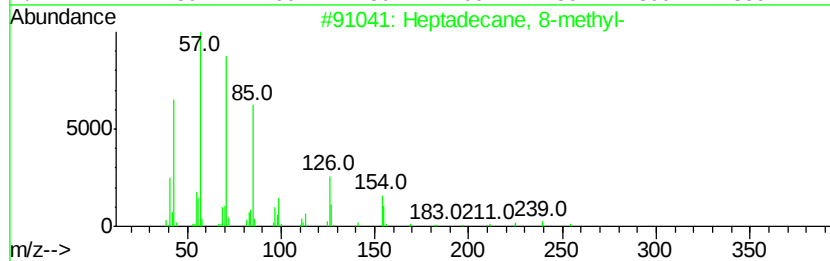
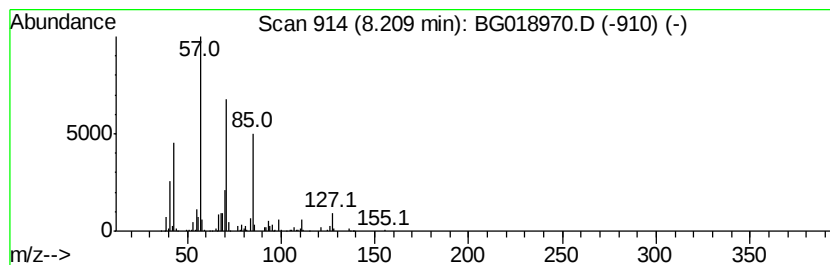
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 8.21 | 19.25 ng/ul | 730274 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptadecane, 8-methyl-   | 254 | C18H38  | 013287-23-5 | 72   |
| 2    |    | Heptacosane              | 380 | C27H56  | 000593-49-7 | 72   |
| 3    |    | Octane, 2,4,6-trimethyl- | 156 | C11H24  | 062016-37-9 | 64   |
| 4    |    | Dodecane, 3-methyl-      | 184 | C13H28  | 017312-57-1 | 64   |
| 5    |    | Decane, 3-methyl-        | 156 | C11H24  | 013151-34-3 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

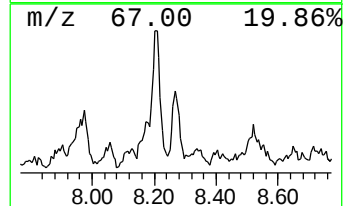
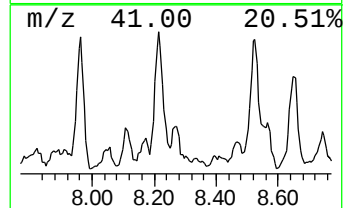
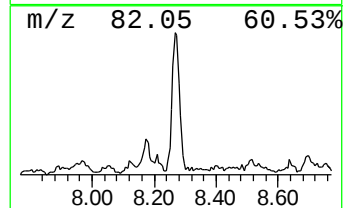
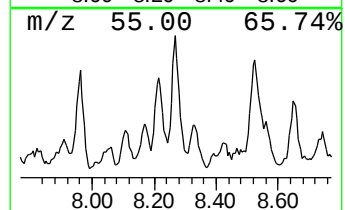
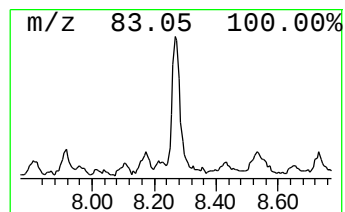
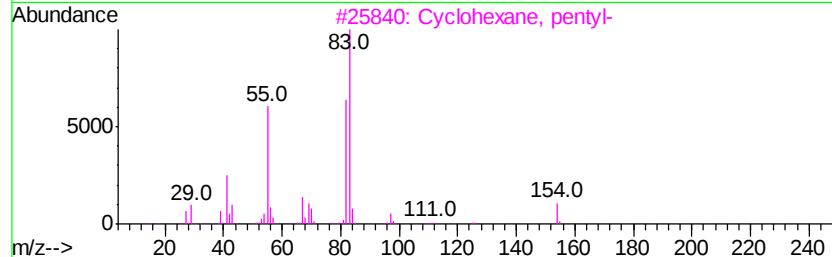
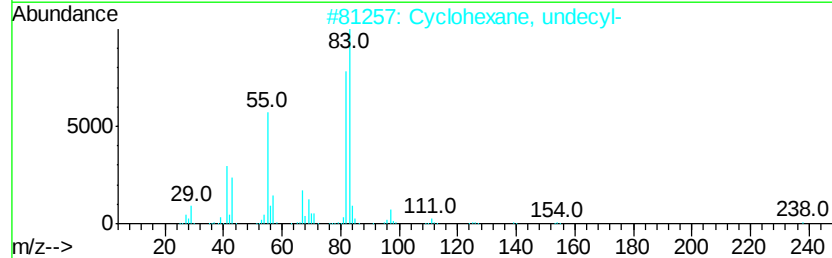
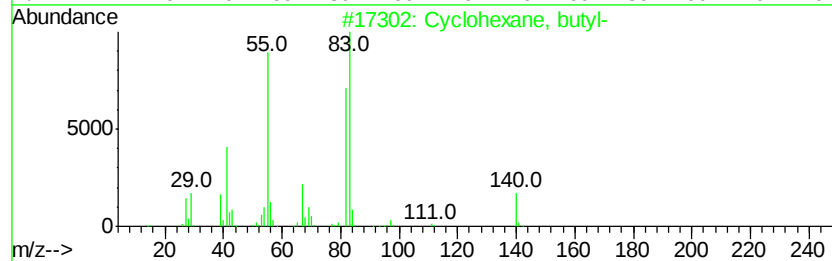
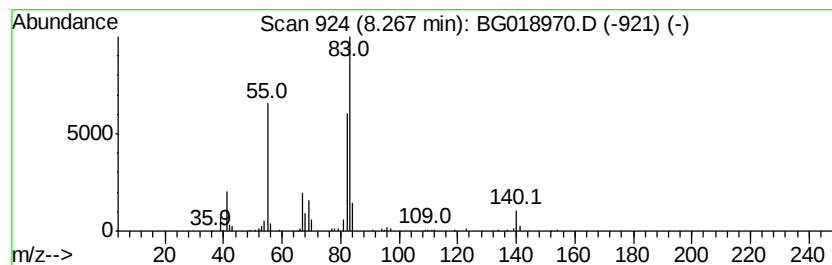
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 (DEL) Alkane: Cyclic8.27 Concentration Rank 53

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.27 | 4.58 ng/ul | 173809 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, butyl-            | 140 | C10H20  | 001678-93-9 | 74   |
| 2    |    | Cyclohexane, undecyl-          | 238 | C17H34  | 054105-66-7 | 72   |
| 3    |    | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 72   |
| 4    |    | Cyclohexane, octyl-            | 196 | C14H28  | 001795-15-9 | 64   |
| 5    |    | Cyclohexane, (2-methylpropyl)- | 140 | C10H20  | 001678-98-4 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

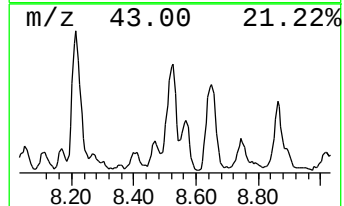
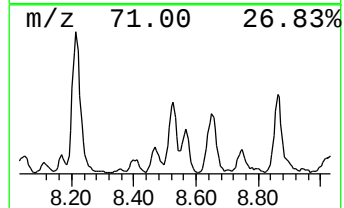
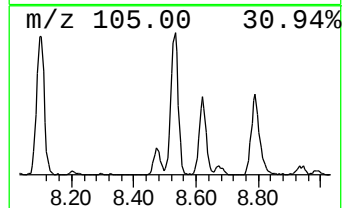
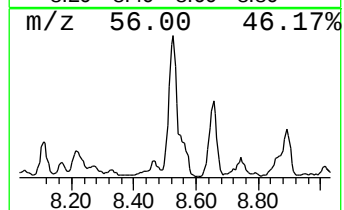
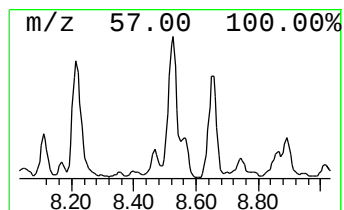
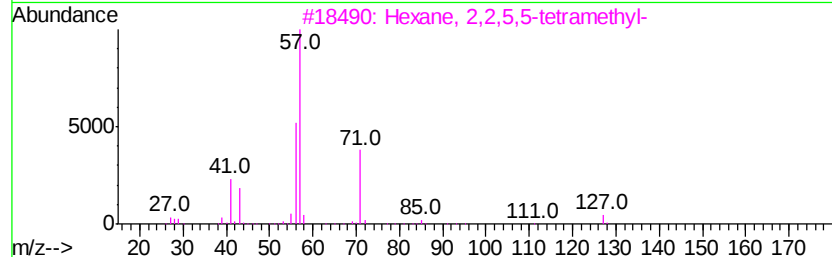
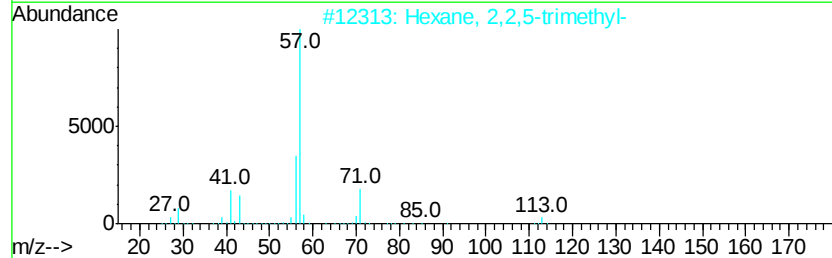
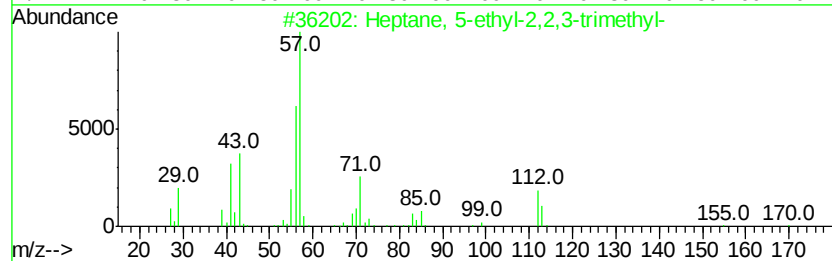
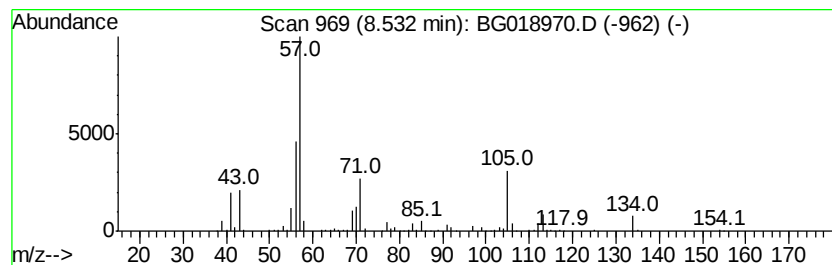
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 8.53 | 17.05 ng/ul | 646849 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptane, 5-ethyl-2,2,3-trimethyl- | 170 | C12H26  | 062199-06-8 | 59   |
| 2    |    | Hexane, 2,2,5-trimethyl-          | 128 | C9H20   | 003522-94-9 | 50   |
| 3    |    | Hexane, 2,2,5,5-tetramethyl-      | 142 | C10H22  | 001071-81-4 | 47   |
| 4    |    | Heptane, 2,2,4,6,6-pentamethyl-   | 170 | C12H26  | 013475-82-6 | 47   |
| 5    |    | Heptane, 2,2,3,5-tetramethyl-     | 156 | C11H24  | 061868-42-6 | 47   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

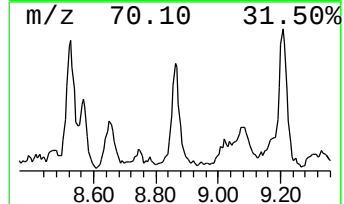
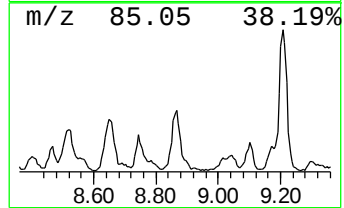
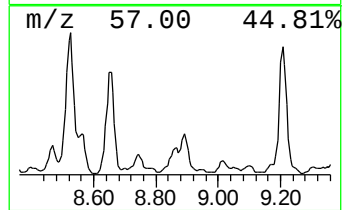
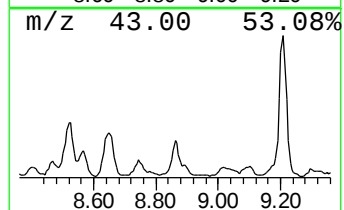
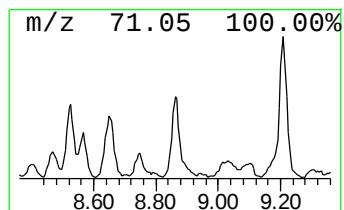
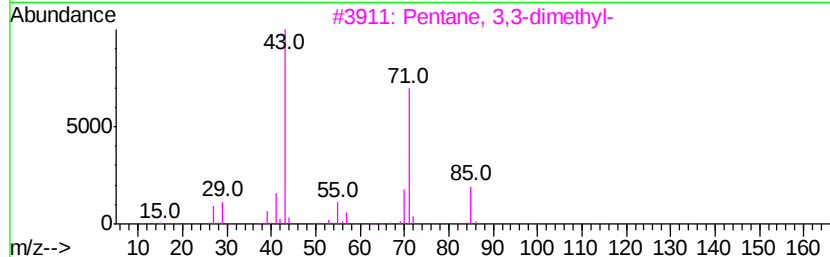
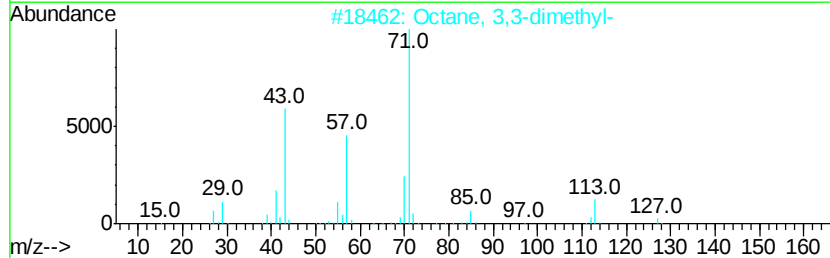
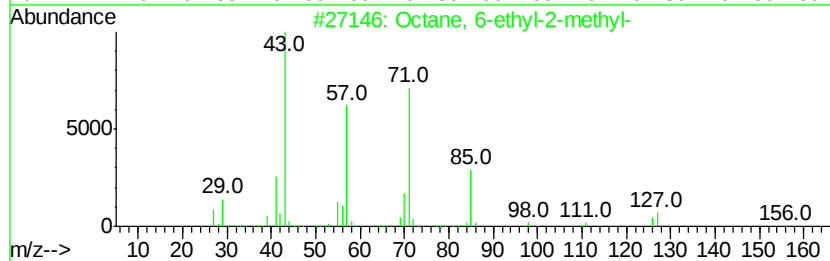
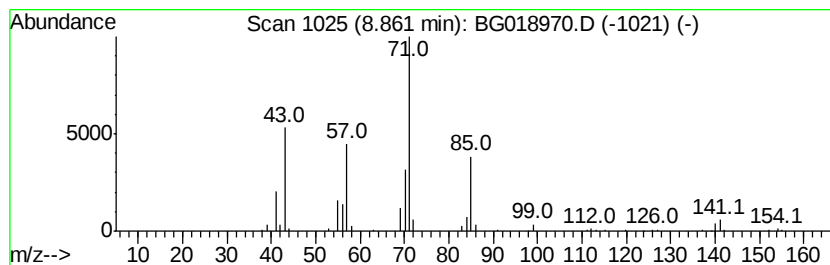
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 55

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.86 | 4.40 ng/ul | 166921 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Octane, 6-ethyl-2-methyl-           | 156 | C11H24   | 062016-19-7  | 56   |
| 2    |    |   | Octane, 3,3-dimethyl-               | 142 | C10H22   | 004110-44-5  | 50   |
| 3    |    |   | Pentane, 3,3-dimethyl-              | 100 | C7H16    | 000562-49-2  | 50   |
| 4    |    |   | Octane, 4-methyl-                   | 128 | C9H20    | 002216-34-4  | 47   |
| 5    |    |   | 1,1-Dimethylpropyl 2-ethylhexanoate | 214 | C13H26O2 | 1000099-99-2 | 47   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

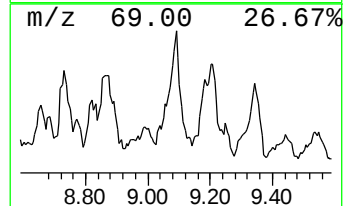
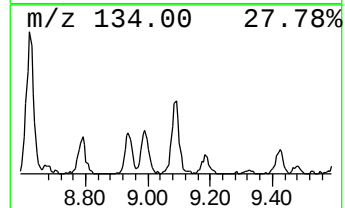
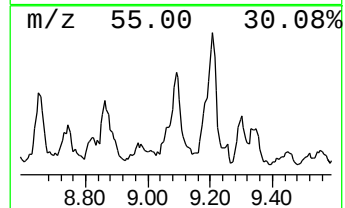
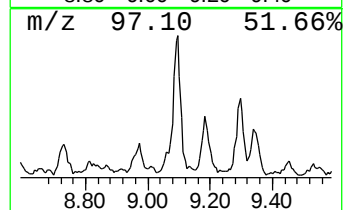
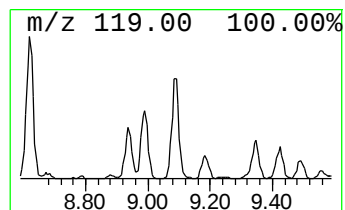
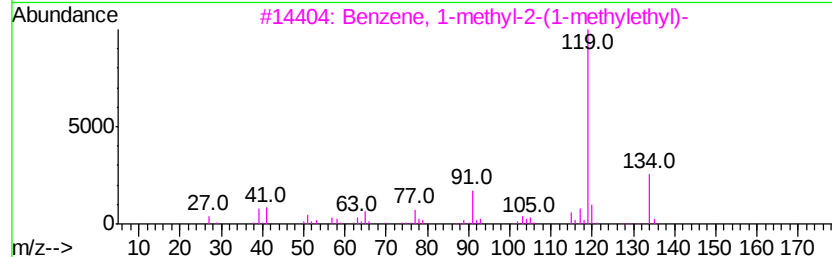
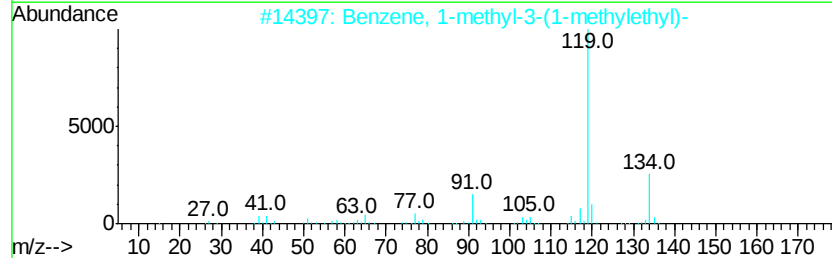
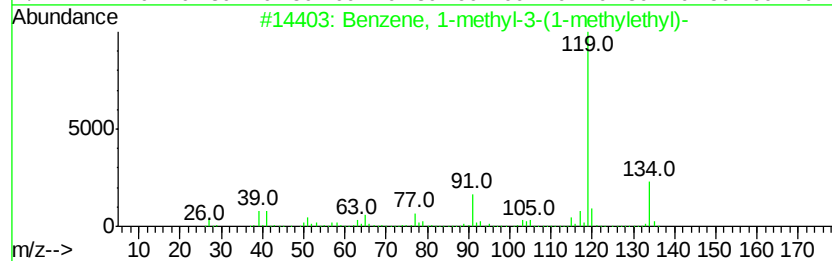
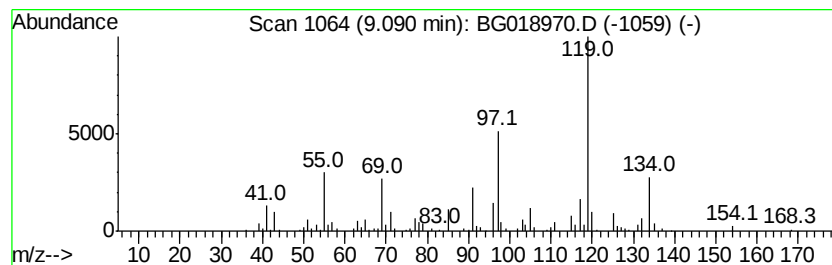
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Benzene, 1-methyl-3-(1-meth... Concentration Rank 37

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.09 | 6.79 ng/ul | 257686 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 90   |
| 2    |    | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 90   |
| 3    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 90   |
| 4    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 90   |
| 5    |    | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

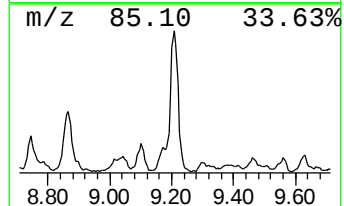
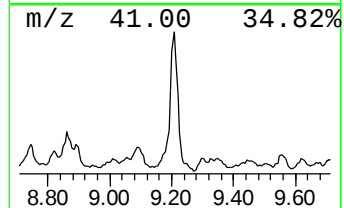
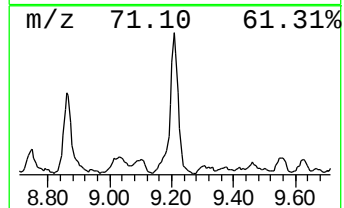
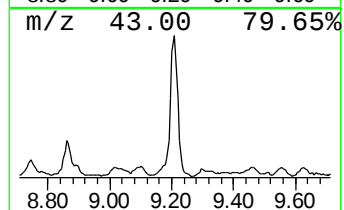
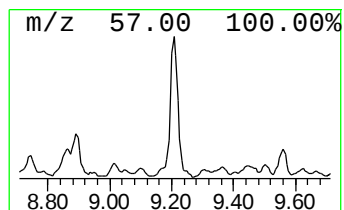
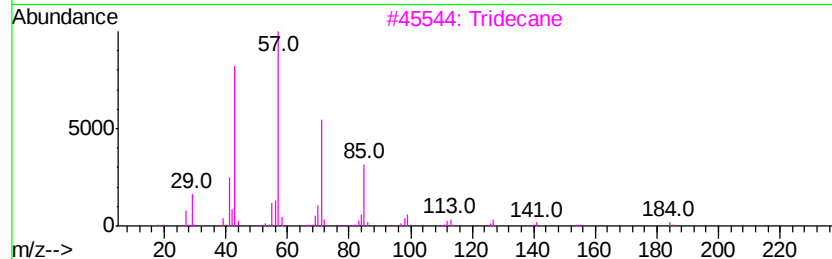
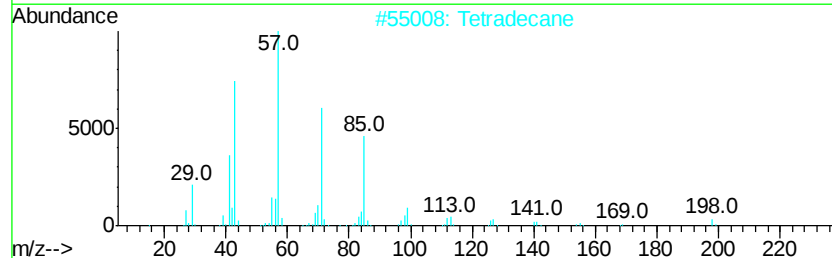
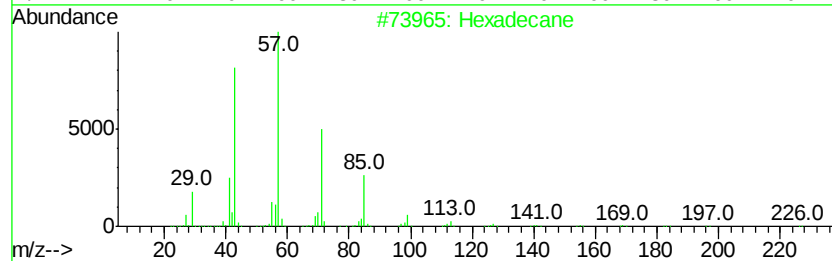
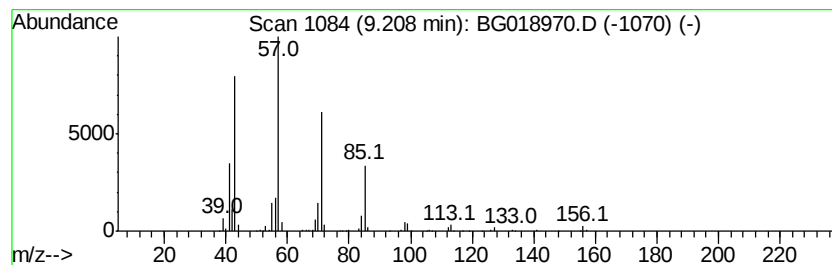
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 9.21 | 26.41 ng/ul | 1001710 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 2    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 83   |
| 3    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 83   |
| 4    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 83   |
| 5    |    | Undecane     | 156 | C11H24  | 001120-21-4 | 76   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

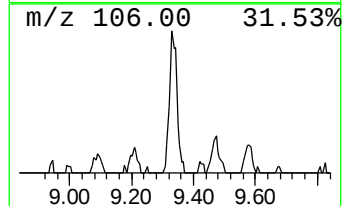
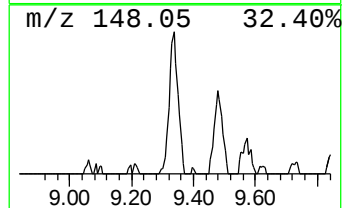
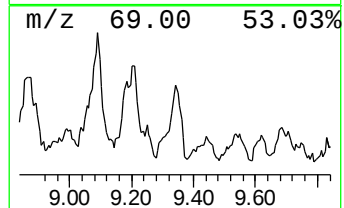
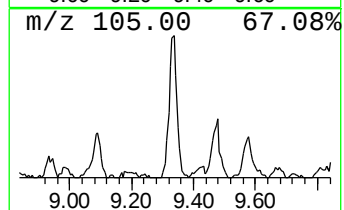
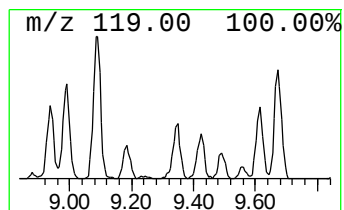
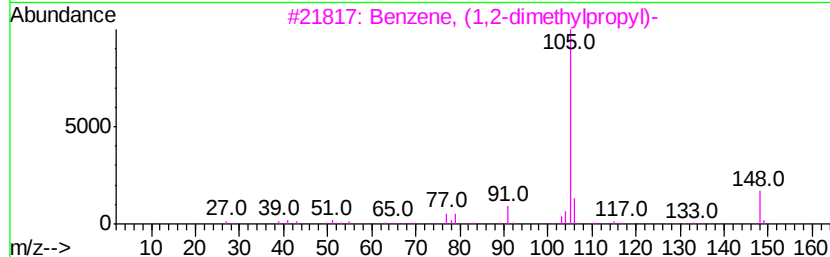
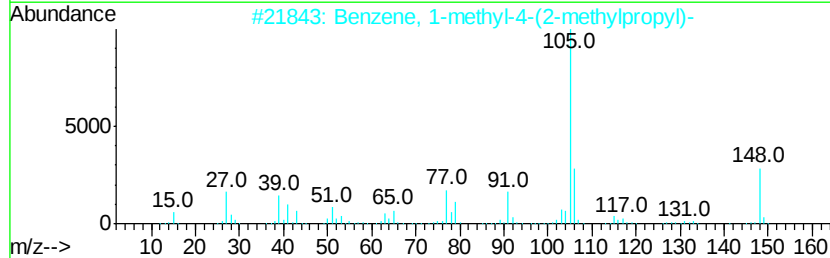
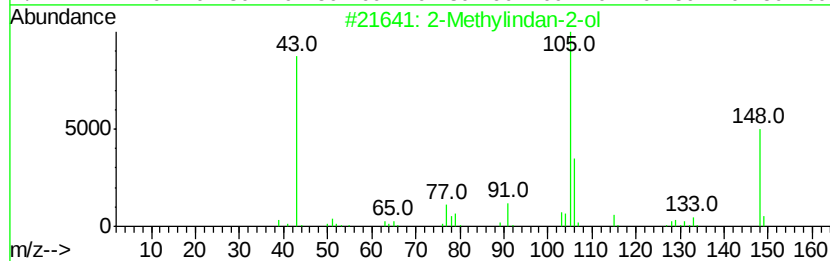
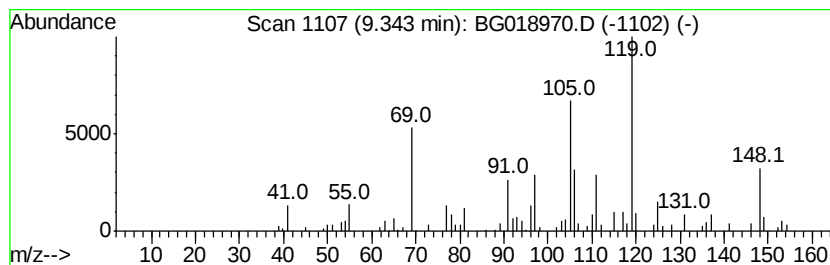
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 unknown-02 Concentration Rank 50

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.34 | 5.08 ng/ul | 192788 | 1,4-Dichlorobenzene-d4 | 7.99 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Methylindan-2-ol                  | 148 | C10H12O  | 033223-84-6 | 41   |
| 2    |    |   | Benzene, 1-methyl-4-(2-methylpro... | 148 | C11H16   | 005161-04-6 | 35   |
| 3    |    |   | Benzene, (1,2-dimethylpropyl)-      | 148 | C11H16   | 004481-30-5 | 22   |
| 4    |    |   | Benzenebutanoic acid, .gamma.-me... | 192 | C12H16O2 | 020881-29-2 | 22   |
| 5    |    |   | Benzene, (1-methylbutyl)-           | 148 | C11H16   | 002719-52-0 | 22   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

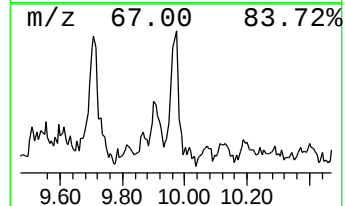
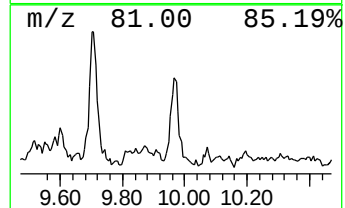
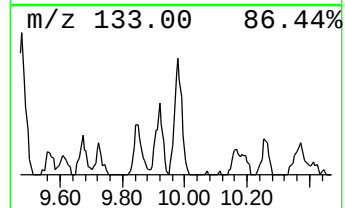
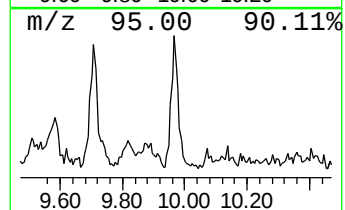
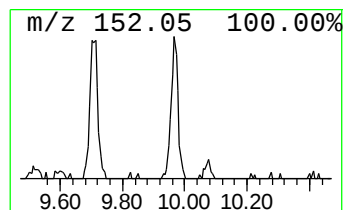
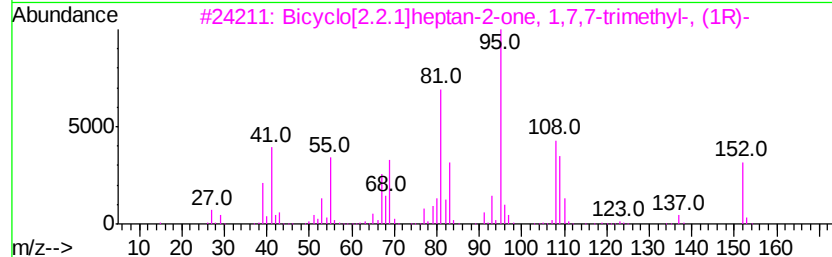
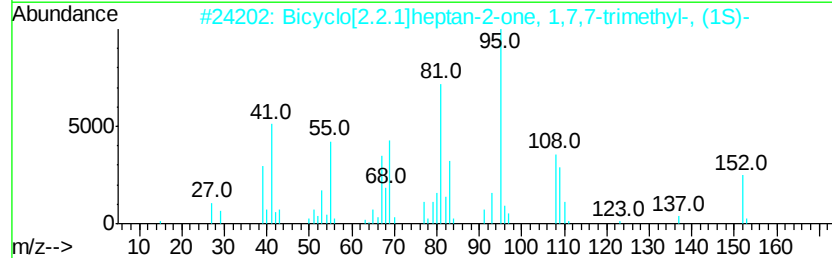
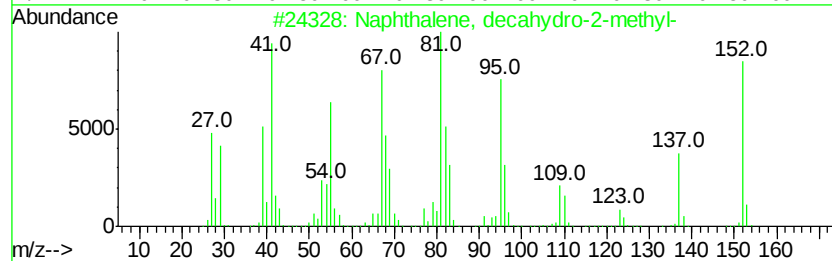
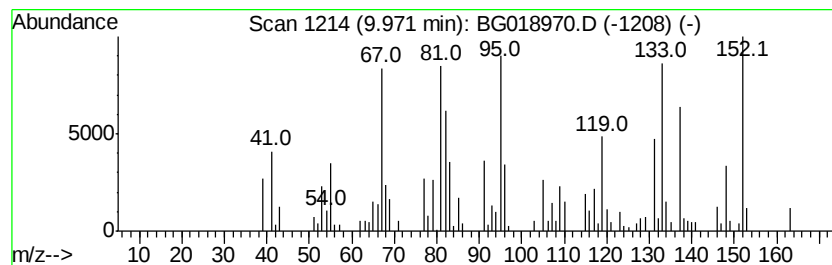
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Naphthalene, decahydro-2-me... Concentration Rank 47

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.97 | 5.44 ng/ul | 147348 | Naphthalene-d8   | 10.79 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Naphthalene, decahydro-2-methyl-    | 152 | C11H20  | 002958-76-1  | 78   |
| 2    |    | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-48-2  | 55   |
| 3    |    | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-49-3  | 55   |
| 4    |    | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-48-2  | 49   |
| 5    |    | cis-Decalin, 2-syn-methyl-          | 152 | C11H20  | 1000155-85-6 | 46   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

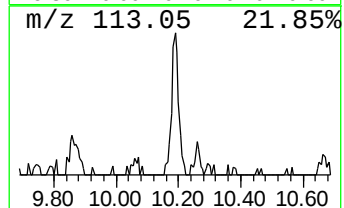
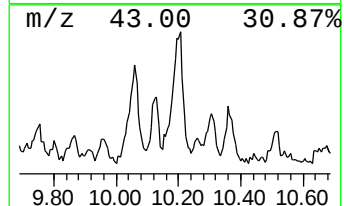
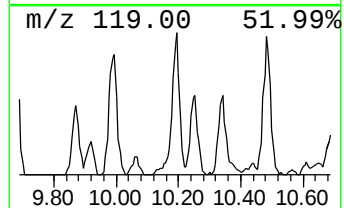
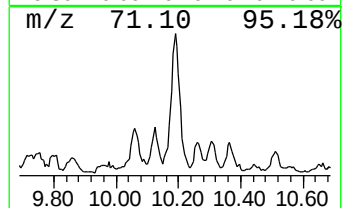
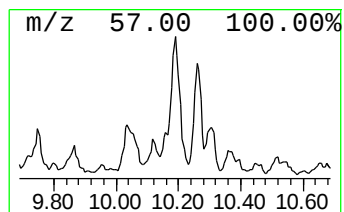
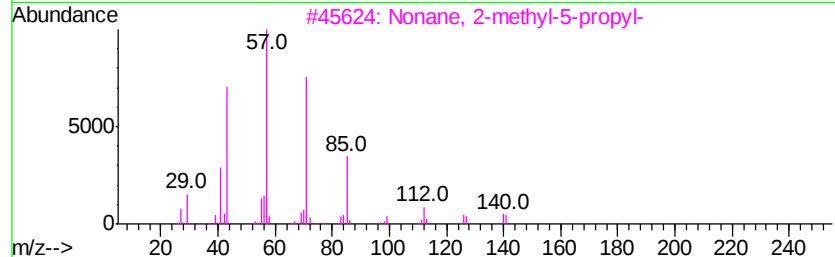
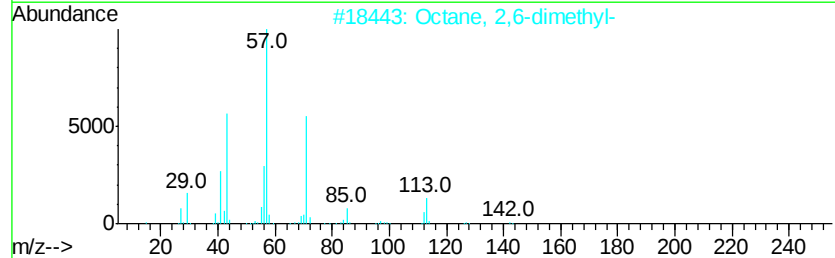
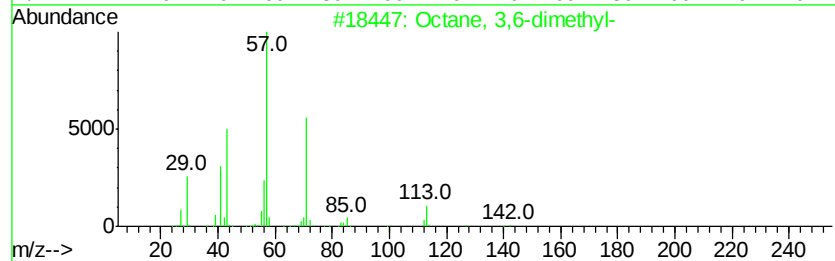
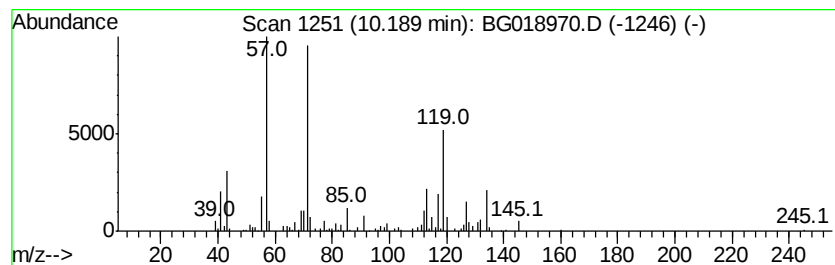
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 34

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.19 | 7.10 ng/ul | 192316 | Naphthalene-d8   | 10.79 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 3,6-dimethyl-          | 142 | C10H22  | 015869-94-0 | 47   |
| 2    |    |   | Octane, 2,6-dimethyl-          | 142 | C10H22  | 002051-30-1 | 38   |
| 3    |    |   | Nonane, 2-methyl-5-propyl-     | 184 | C13H28  | 031081-17-1 | 35   |
| 4    |    |   | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 11   |
| 5    |    |   | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14  | 002870-04-4 | 11   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

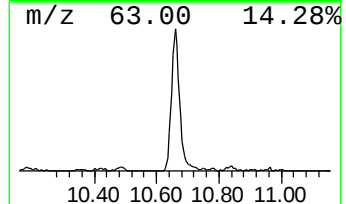
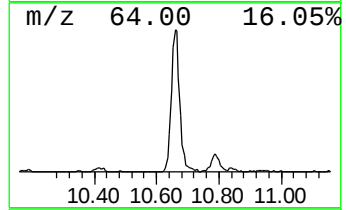
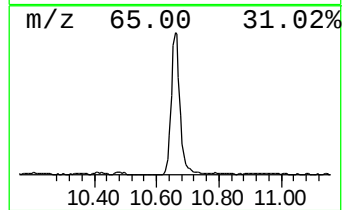
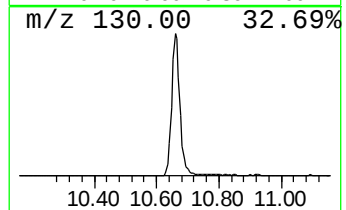
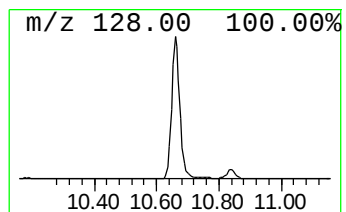
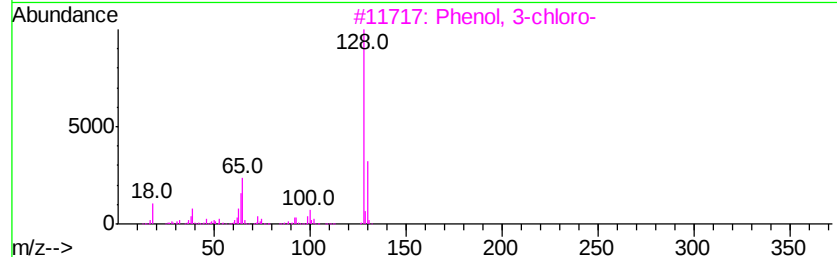
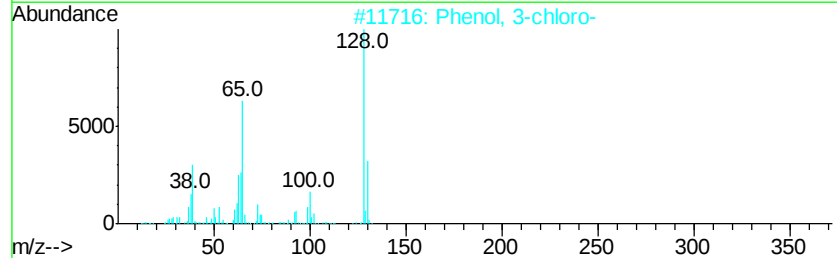
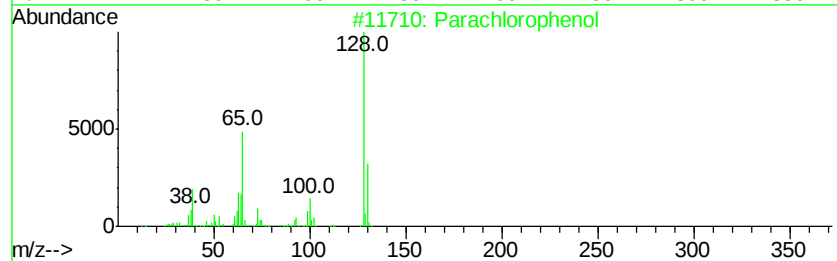
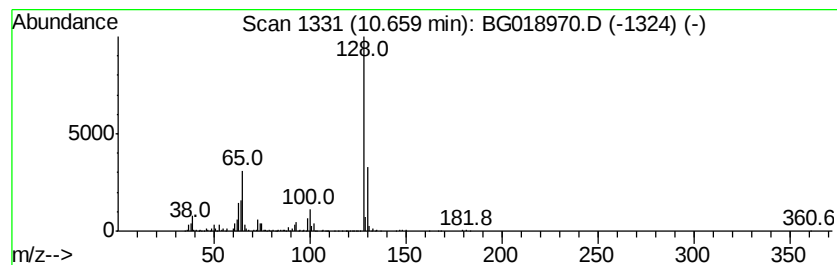
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Parachlorophenol Concentration Rank 3

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 10.66 | 35.17 ng/ul | 952422 | Naphthalene-d8   | 10.79 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Parachlorophenol  | 128 | C6H5ClO | 000106-48-9 | 95   |
| 2    |    | Phenol, 3-chloro- | 128 | C6H5ClO | 000108-43-0 | 95   |
| 3    |    | Phenol, 3-chloro- | 128 | C6H5ClO | 000108-43-0 | 95   |
| 4    |    | Parachlorophenol  | 128 | C6H5ClO | 000106-48-9 | 94   |
| 5    |    | Phenol, 3-chloro- | 128 | C6H5ClO | 000108-43-0 | 94   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

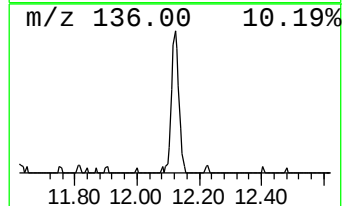
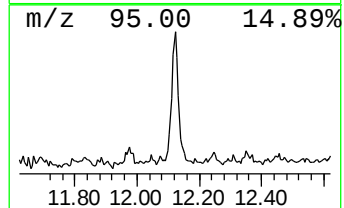
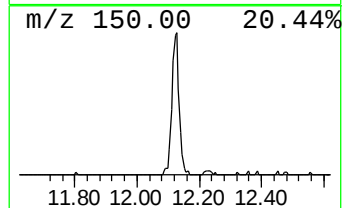
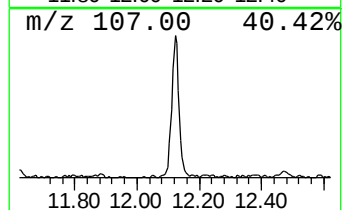
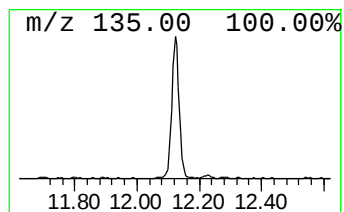
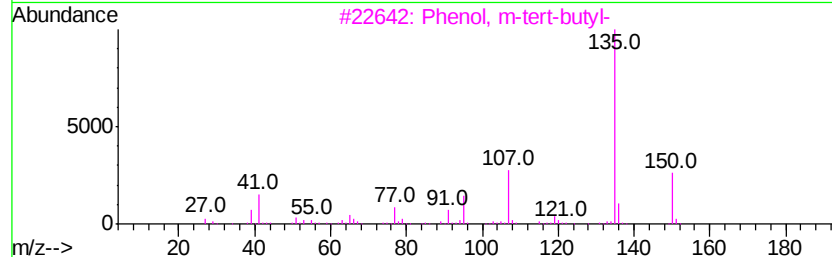
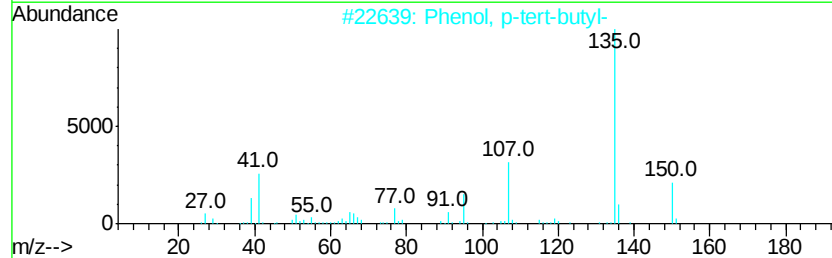
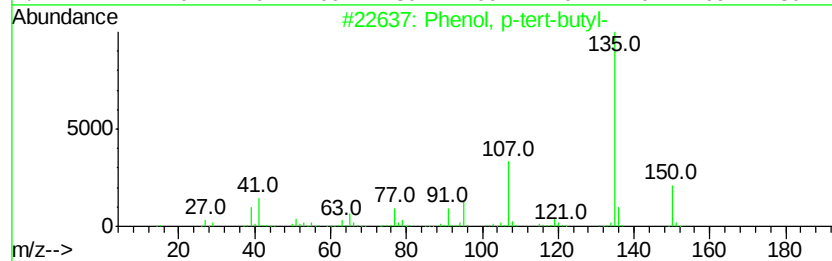
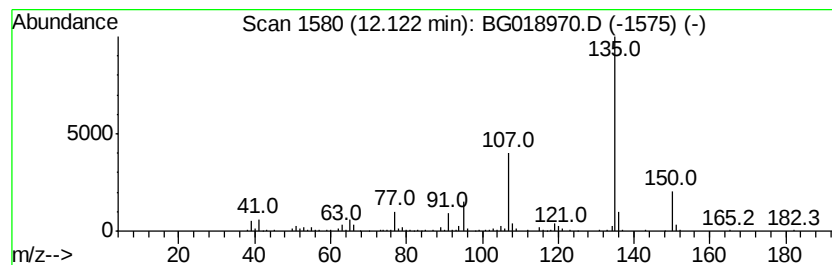
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Phenol, p-tert-butyl- Concentration Rank 33

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.12 | 7.37 ng/ul | 199577 | Naphthalene-d8   | 10.79 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 96   |
| 2    |    | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 94   |
| 3    |    | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 93   |
| 4    |    | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 91   |
| 5    |    | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 91   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

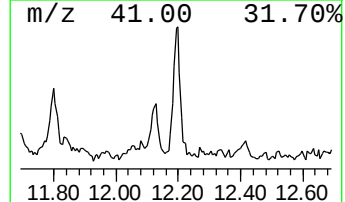
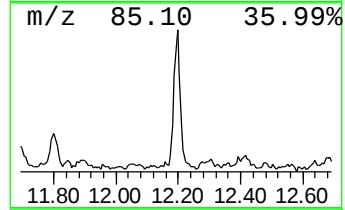
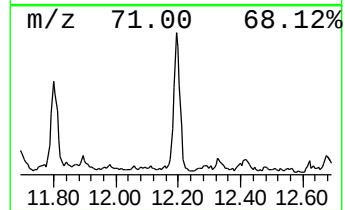
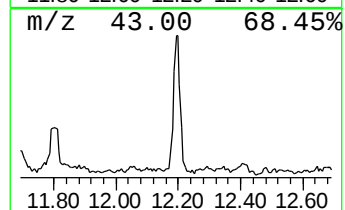
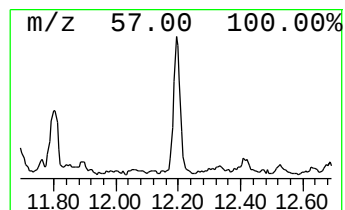
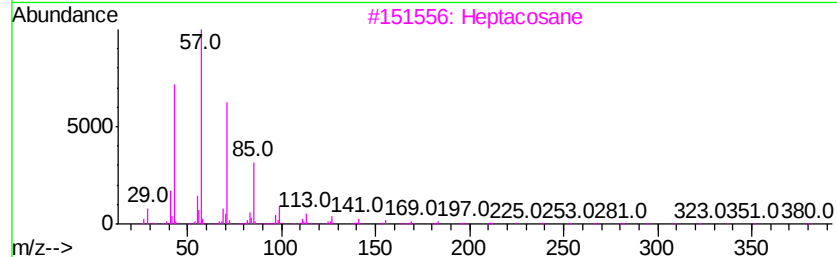
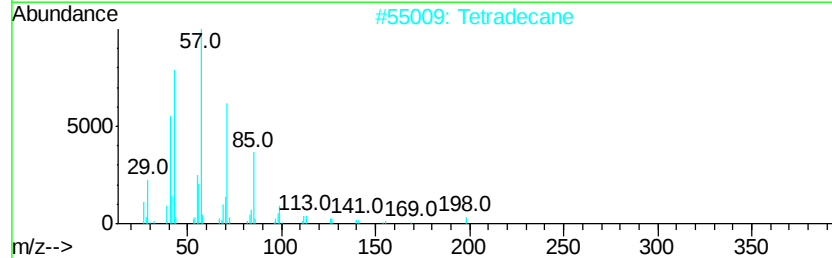
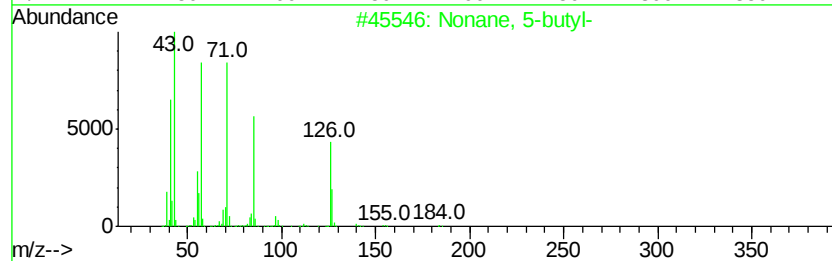
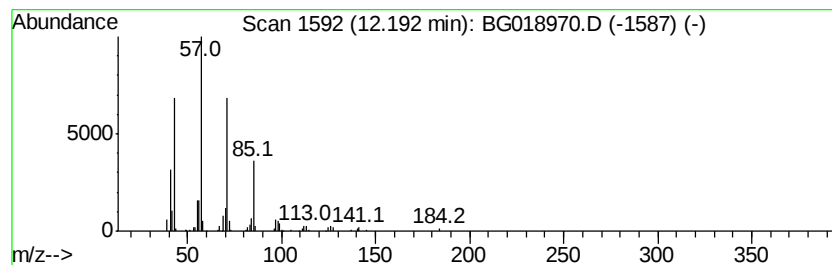
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.19 | 6.27 ng/ul | 169942 | Napthalene-d8    | 10.79 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 5-butyl-     | 184 | C13H28  | 017312-63-9 | 87   |
| 2    |    | Tetradecane          | 198 | C14H30  | 000629-59-4 | 86   |
| 3    |    | Heptacosane          | 380 | C27H56  | 000593-49-7 | 78   |
| 4    |    | Eicosane, 10-methyl- | 296 | C21H44  | 054833-23-7 | 78   |
| 5    |    | Nonadecane           | 268 | C19H40  | 000629-92-5 | 78   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

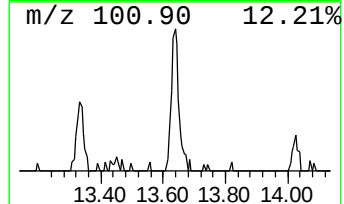
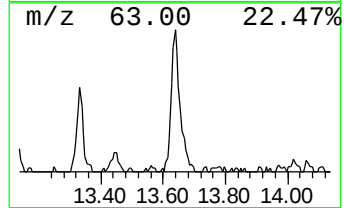
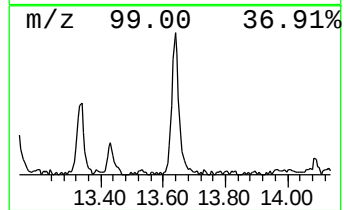
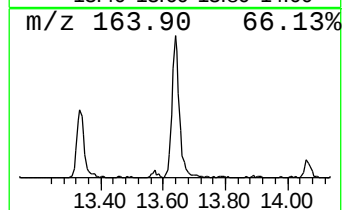
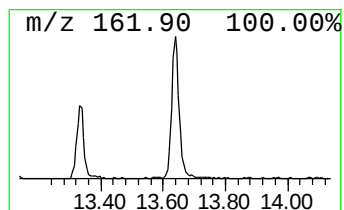
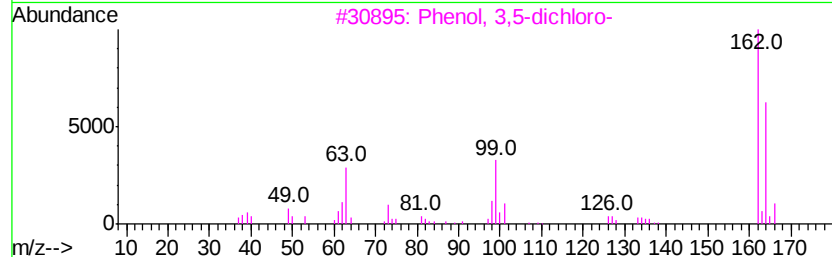
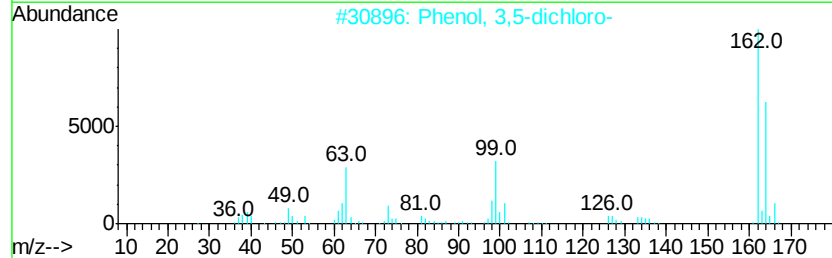
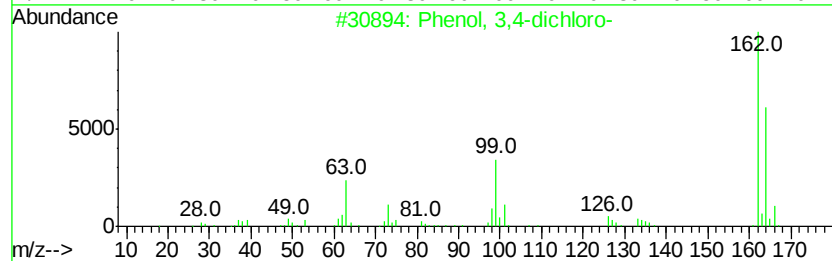
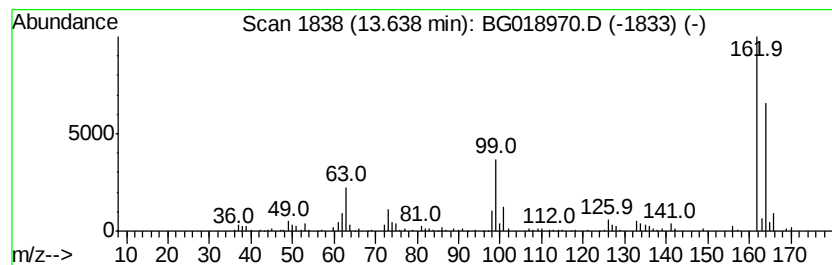
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Phenol, 3,4-dichloro- Concentration Rank 63

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.64 | 3.38 ng/ul | 174155 | Acenaphthene-d10 | 14.61 |

| Hit# | of                    | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|-----------------------|---|--------------|-----|----------|-------------|------|
| 1    | Phenol, 3,4-dichloro- |   |              | 162 | C6H4Cl2O | 000095-77-2 | 98   |
| 2    | Phenol, 3,5-dichloro- |   |              | 162 | C6H4Cl2O | 000591-35-5 | 96   |
| 3    | Phenol, 3,5-dichloro- |   |              | 162 | C6H4Cl2O | 000591-35-5 | 96   |
| 4    | Phenol, 3,4-dichloro- |   |              | 162 | C6H4Cl2O | 000095-77-2 | 95   |
| 5    | Phenol, 3,5-dichloro- |   |              | 162 | C6H4Cl2O | 000591-35-5 | 91   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

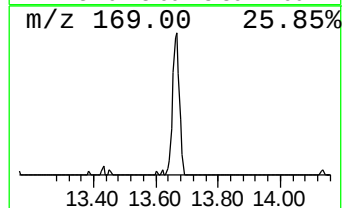
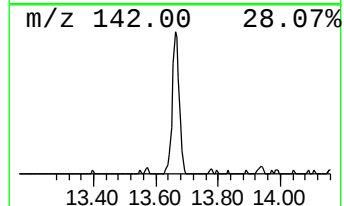
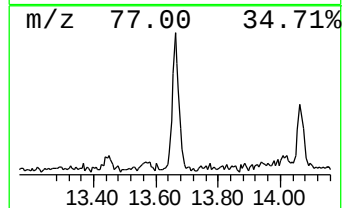
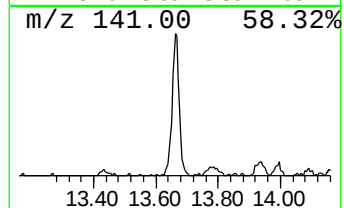
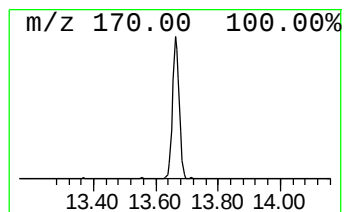
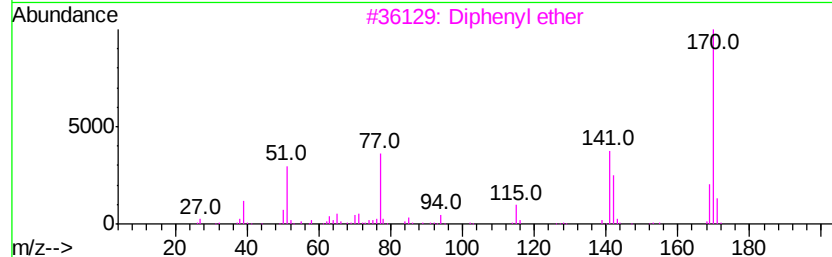
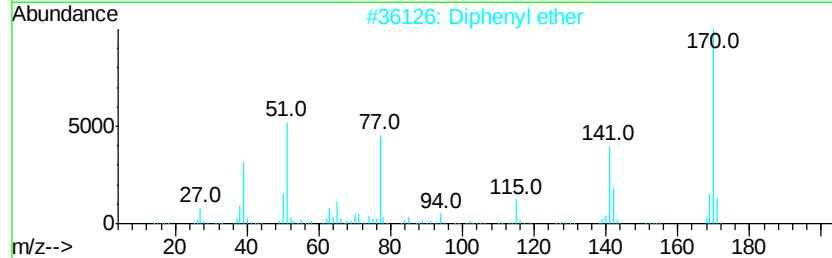
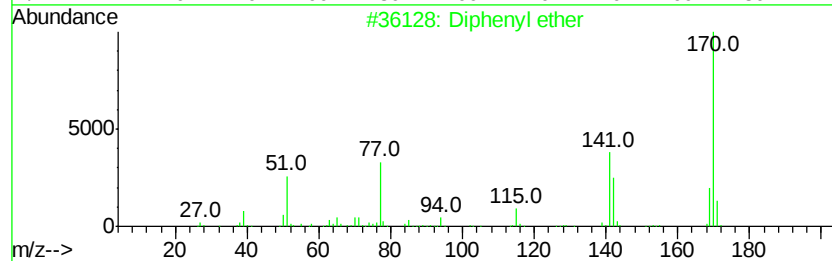
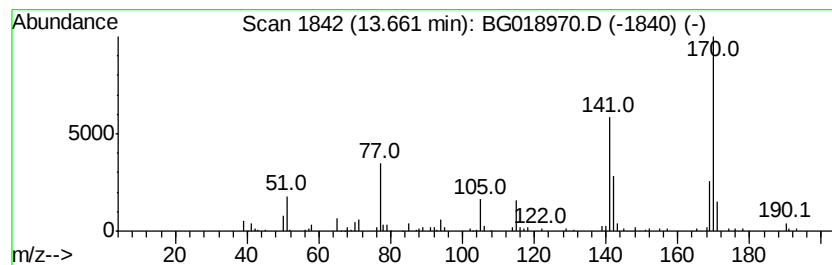
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 Diphenyl ether Concentration Rank 60

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.66 | 3.47 ng/ul | 178469 | Acenaphthene-d10 | 14.61 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Diphenyl ether              | 170 | C12H10O | 000101-84-8 | 93   |
| 2    |    | Diphenyl ether              | 170 | C12H10O | 000101-84-8 | 87   |
| 3    |    | Diphenyl ether              | 170 | C12H10O | 000101-84-8 | 87   |
| 4    |    | Diphenyl ether              | 170 | C12H10O | 000101-84-8 | 81   |
| 5    |    | 1,2-Dehydro-as-hydrindacene | 170 | C13H14  | 096508-75-7 | 53   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

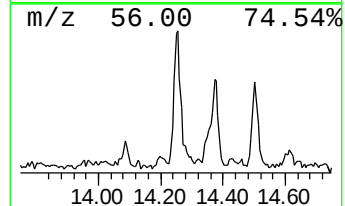
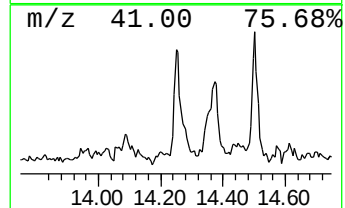
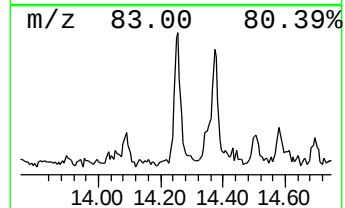
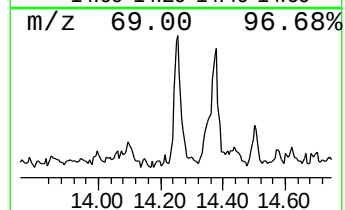
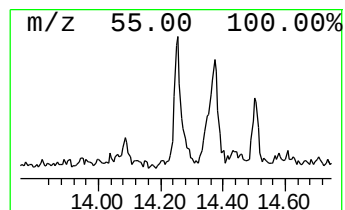
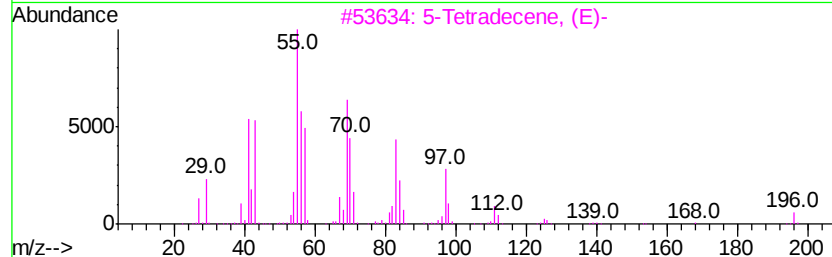
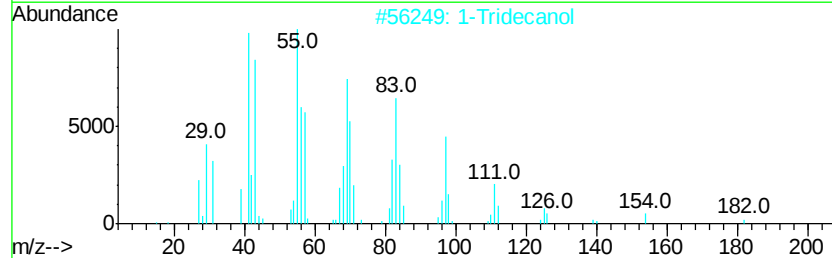
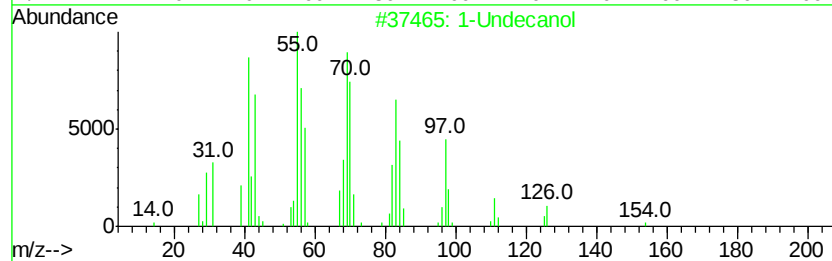
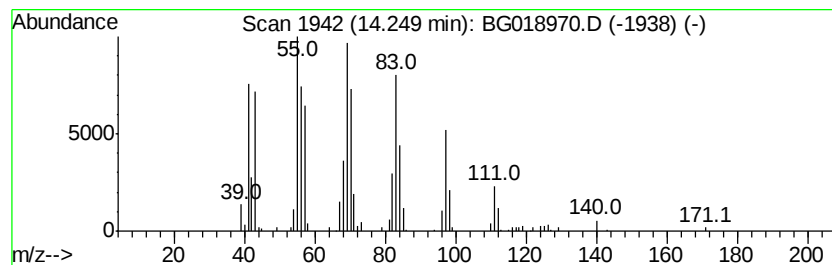
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 1-Undecanol Concentration Rank 66

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.25 | 3.14 ng/ul | 161716 | Acenaphthene-d10 | 14.61 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Undecanol         | 172 | C11H24O | 000112-42-5 | 91   |
| 2    |    |   | 1-Tridecanol        | 200 | C13H28O | 000112-70-9 | 91   |
| 3    |    |   | 5-Tetradecene, (E)- | 196 | C14H28  | 041446-66-6 | 91   |
| 4    |    |   | Cyclododecane       | 140 | C10H20  | 000293-96-9 | 90   |
| 5    |    |   | Cyclododecane       | 168 | C12H24  | 000294-62-2 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

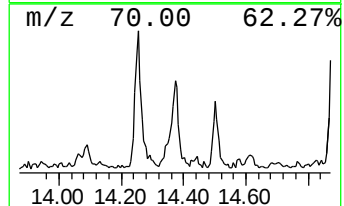
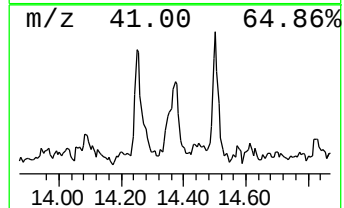
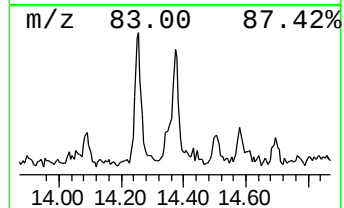
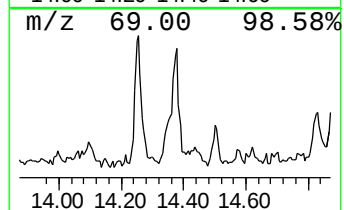
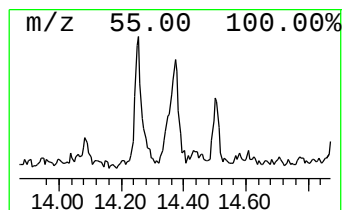
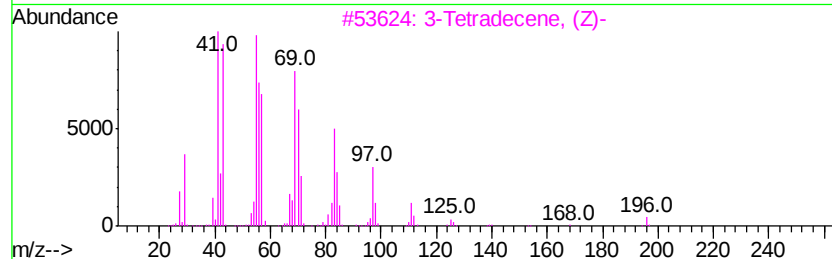
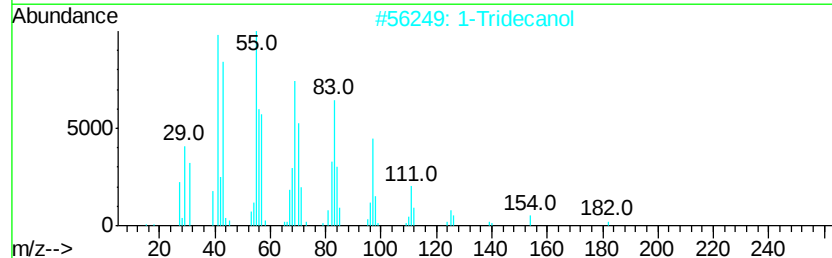
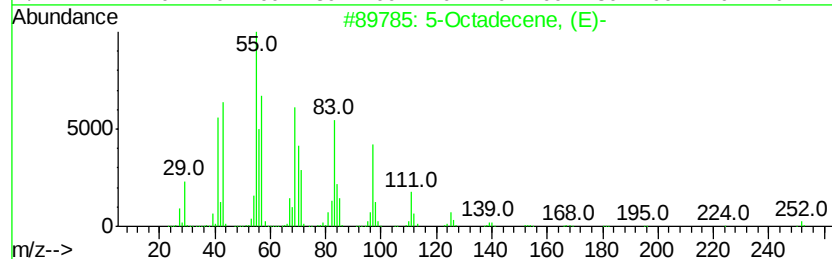
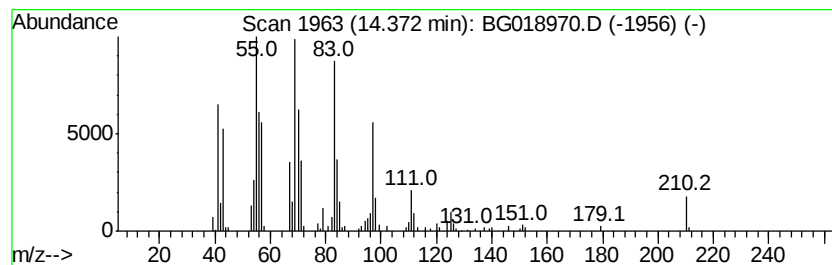
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 (DEL) Alkane: Cyclic14.37 Concentration Rank 57

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.37 | 4.02 ng/ul | 207050 | Acenaphthene-d10 | 14.61 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | 5-Octadecene, (E)-  | 252 | C18H36  | 007206-21-5 | 91   |
| 2    |    | 1-Tridecanol        | 200 | C13H28O | 000112-70-9 | 90   |
| 3    |    | 3-Tetradecene, (Z)- | 196 | C14H28  | 041446-67-7 | 90   |
| 4    |    | 5-Tetradecene, (E)- | 196 | C14H28  | 041446-66-6 | 90   |
| 5    |    | 1-Pentadecanol      | 228 | C15H32O | 000629-76-5 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

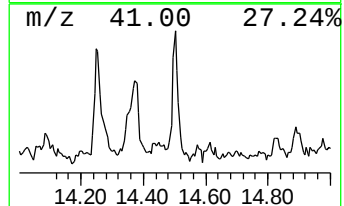
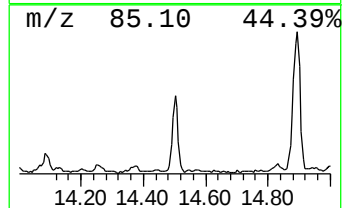
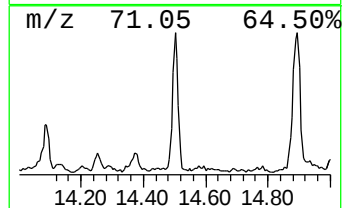
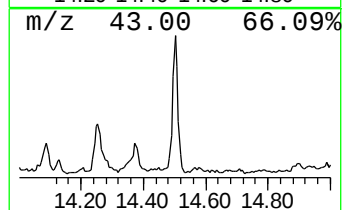
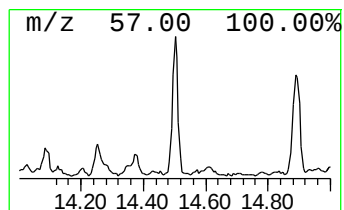
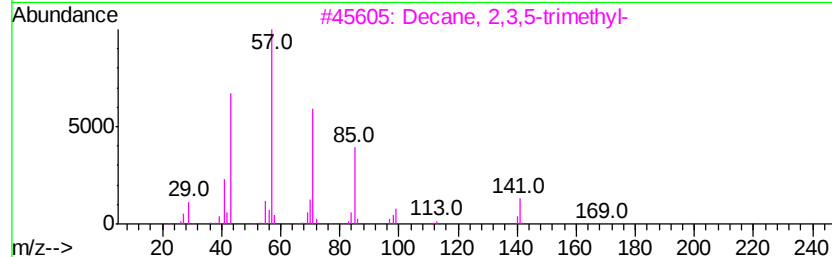
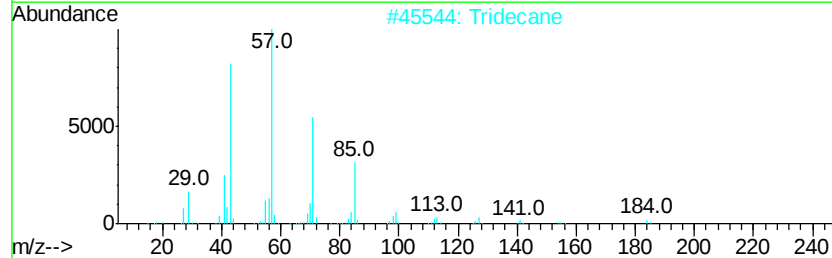
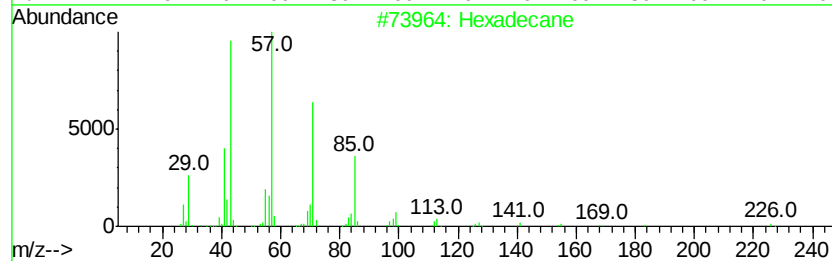
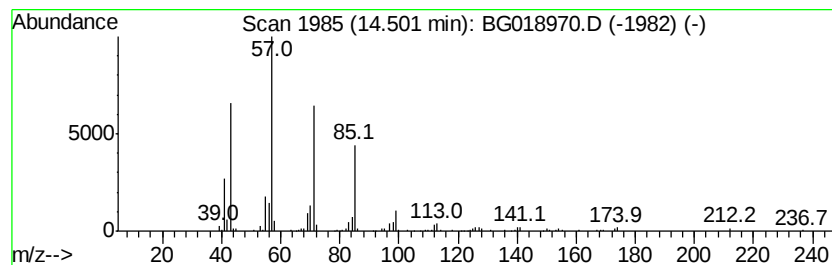
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 62

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.50 | 3.44 ng/ul | 176861 | Acenaphthene-d10 | 14.61 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane               | 226 | C16H34  | 000544-76-3 | 90   |
| 2    |    |   | Tridecane                | 184 | C13H28  | 000629-50-5 | 83   |
| 3    |    |   | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 83   |
| 4    |    |   | Eicosane                 | 282 | C20H42  | 000112-95-8 | 80   |
| 5    |    |   | Hexadecane               | 226 | C16H34  | 000544-76-3 | 78   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

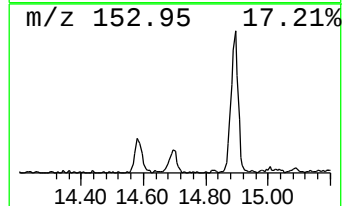
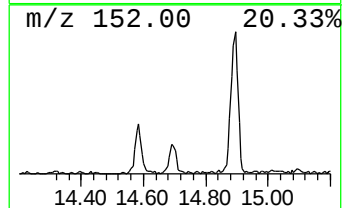
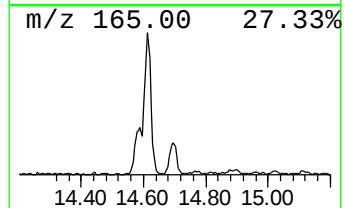
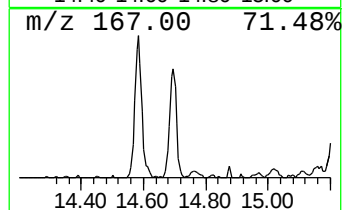
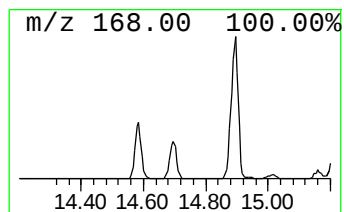
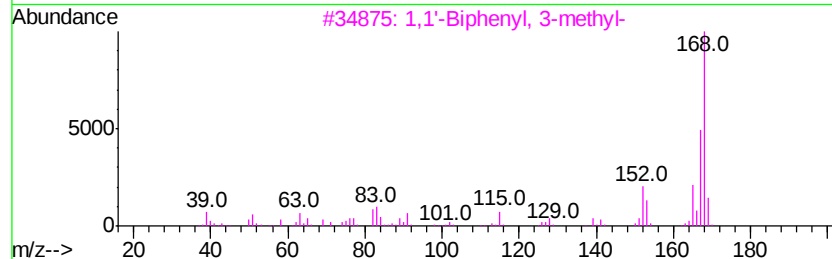
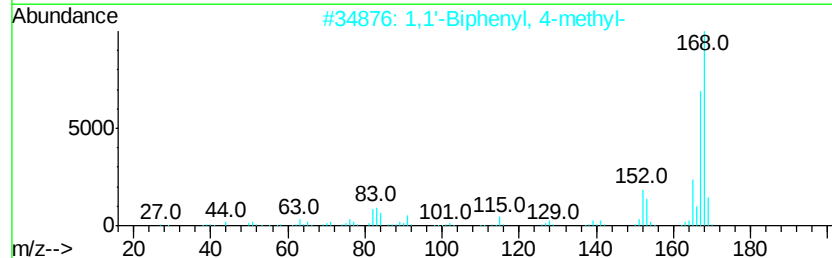
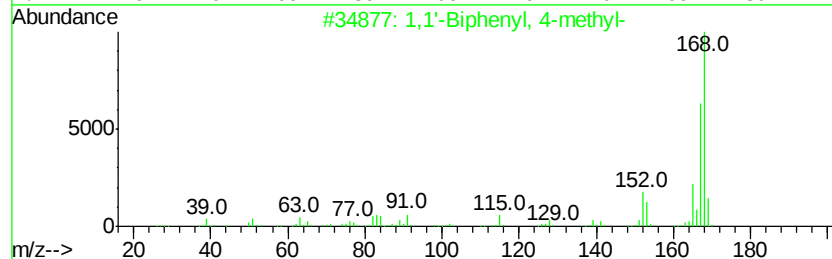
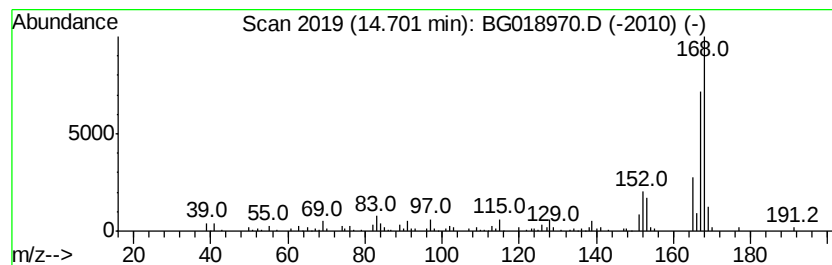
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 1,1'-Biphenyl, 4-methyl- Concentration Rank 71

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.70 | 2.72 ng/ul | 139796 | Acenaphthene-d10 | 14.61 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | 1,1'-Biphenyl, 4-methyl- | 168 | C13H12  | 000644-08-6 | 93   |
| 2    |    |   | 1,1'-Biphenyl, 4-methyl- | 168 | C13H12  | 000644-08-6 | 87   |
| 3    |    |   | 1,1'-Biphenyl, 3-methyl- | 168 | C13H12  | 000643-93-6 | 86   |
| 4    |    |   | 1,1'-Biphenyl, 3-methyl- | 168 | C13H12  | 000643-93-6 | 83   |
| 5    |    |   | Diphenylmethane          | 168 | C13H12  | 000101-81-5 | 83   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

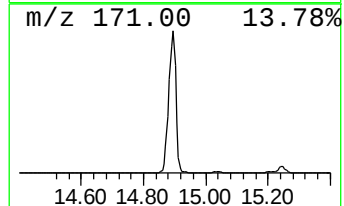
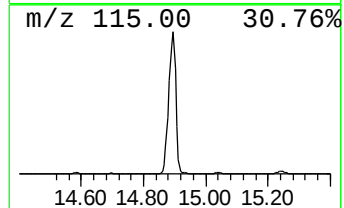
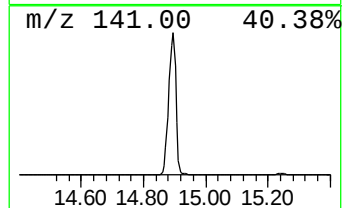
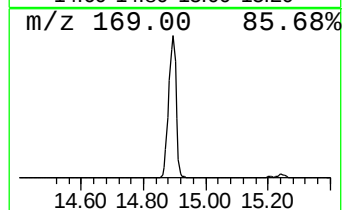
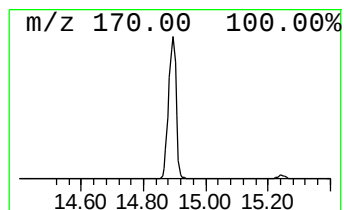
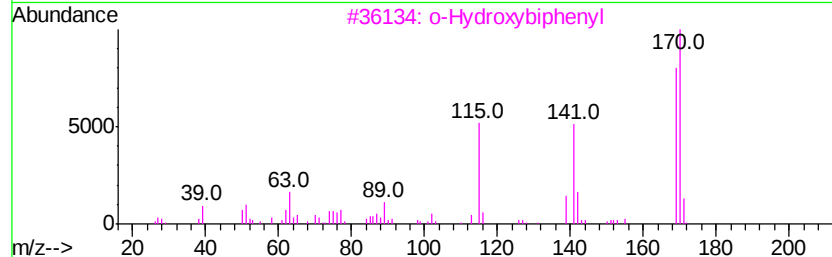
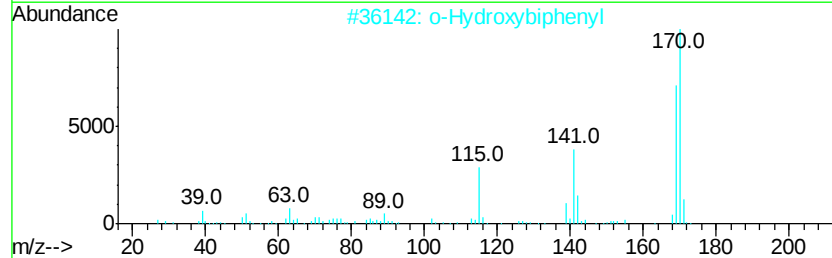
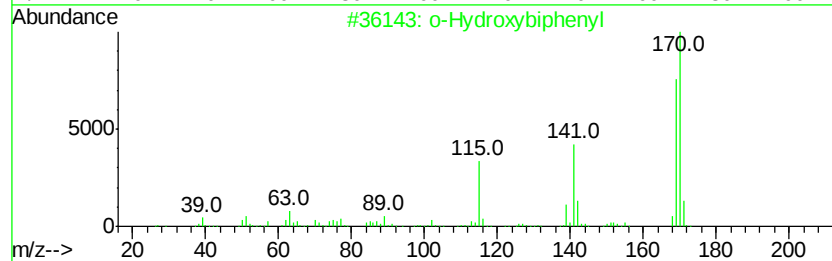
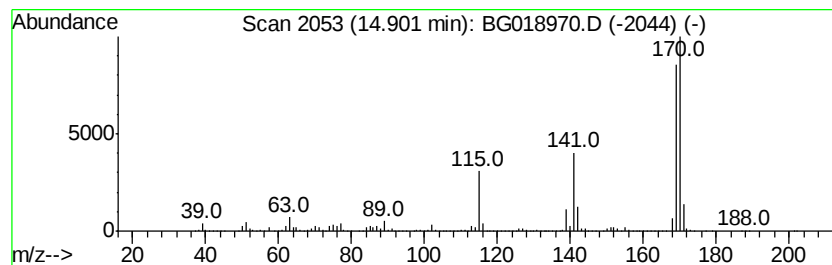
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 o-Hydroxybiphenyl Concentration Rank 1

| R.T.  | EstConc      | Area    | Relative to ISTD | R.T.  |
|-------|--------------|---------|------------------|-------|
| 14.90 | 124.14 ng/ul | 6388840 | Acenaphthene-d10 | 14.61 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | o-Hydroxybiphenyl | 170 | C12H10O | 000090-43-7 | 97   |
| 2    |    | o-Hydroxybiphenyl | 170 | C12H10O | 000090-43-7 | 97   |
| 3    |    | o-Hydroxybiphenyl | 170 | C12H10O | 000090-43-7 | 96   |
| 4    |    | o-Hydroxybiphenyl | 170 | C12H10O | 000090-43-7 | 90   |
| 5    |    | p-Hydroxybiphenyl | 170 | C12H10O | 000092-69-3 | 62   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

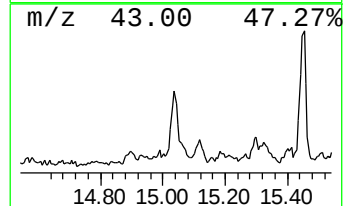
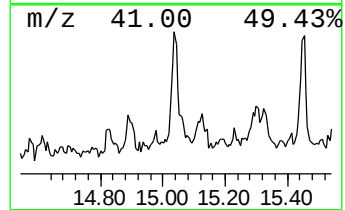
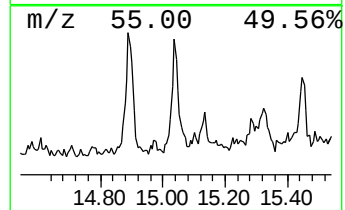
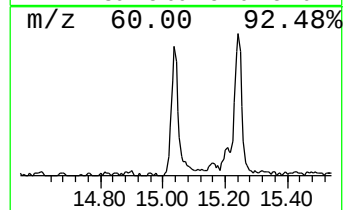
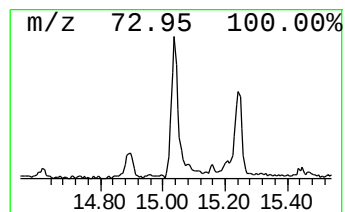
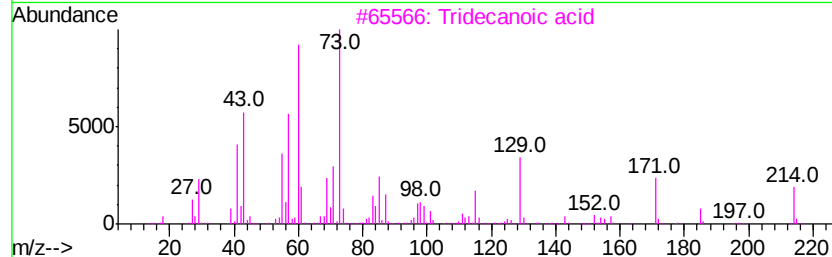
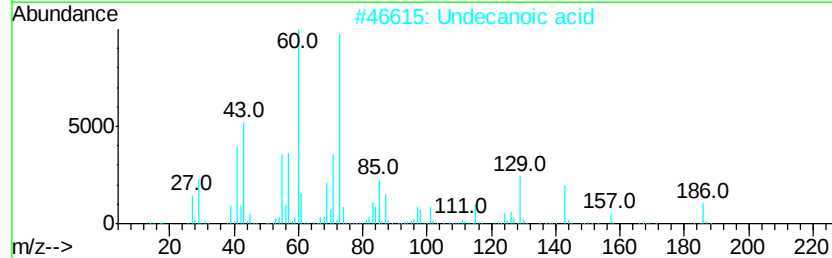
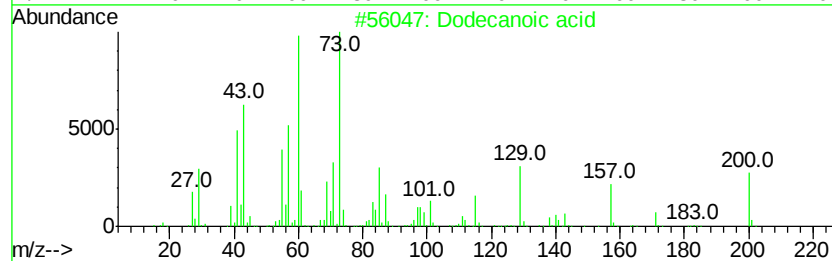
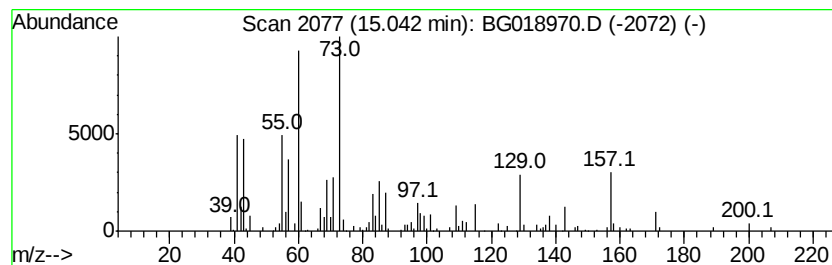
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 Dodecanoic acid Concentration Rank 61

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.04 | 3.45 ng/ul | 177354 | Acenaphthene-d10 | 14.61 |

| Hit# | of | Tentative ID     | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------|-----|----------|-------------|------|
| 1    | 5  | Dodecanoic acid  | 200 | C12H24O2 | 000143-07-7 | 90   |
| 2    |    | Undecanoic acid  | 186 | C11H22O2 | 000112-37-8 | 72   |
| 3    |    | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 64   |
| 4    |    | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 64   |
| 5    |    | n-Decanoic acid  | 172 | C10H20O2 | 000334-48-5 | 59   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

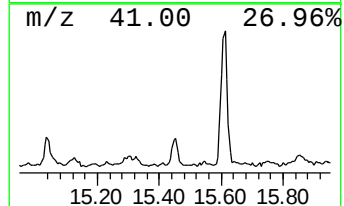
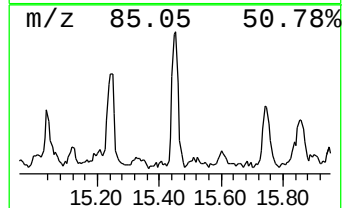
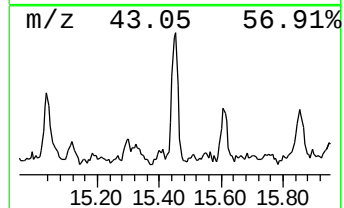
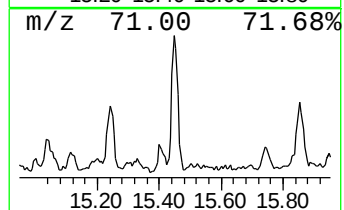
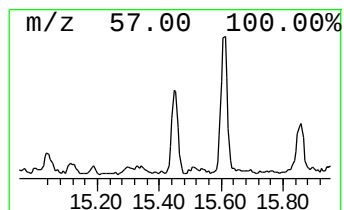
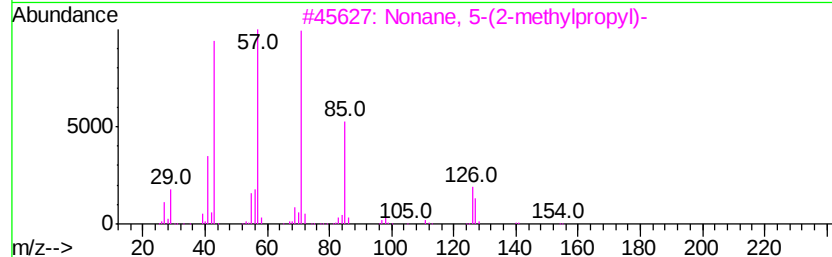
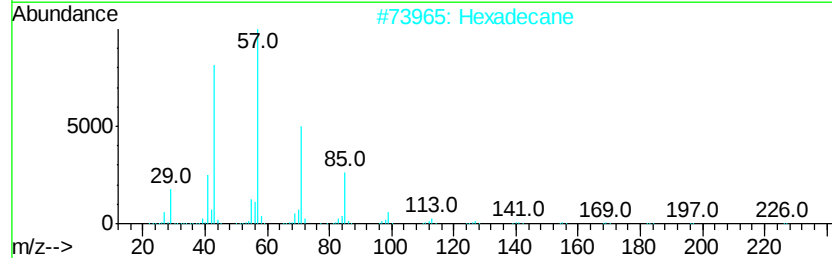
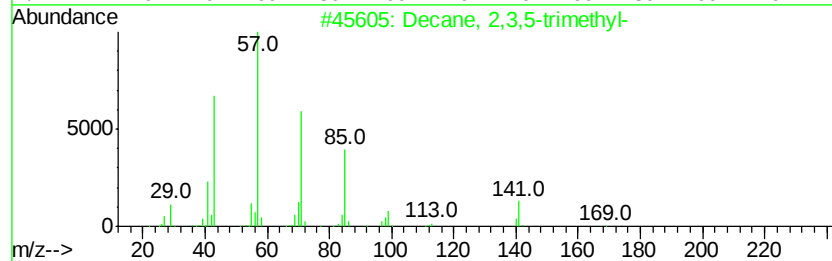
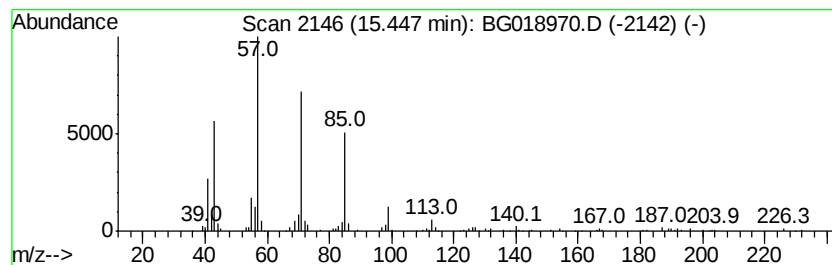
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 70

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.45 | 2.86 ng/ul | 147098 | Acenaphthene-d10 | 14.61 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,3,5-trimethyl-    | 184 | C13H28  | 062238-11-3 | 72   |
| 2    |    |   | Hexadecane                  | 226 | C16H34  | 000544-76-3 | 64   |
| 3    |    |   | Nonane, 5-(2-methylpropyl)- | 184 | C13H28  | 062185-53-9 | 64   |
| 4    |    |   | Tridecane, 2-methyl-        | 198 | C14H30  | 001560-96-9 | 59   |
| 5    |    |   | Hexadecane                  | 226 | C16H34  | 000544-76-3 | 58   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

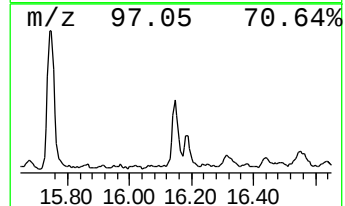
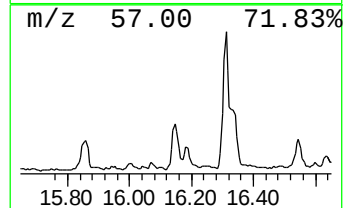
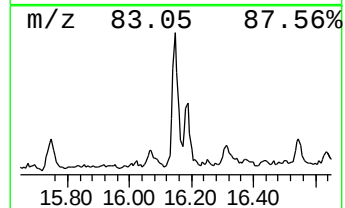
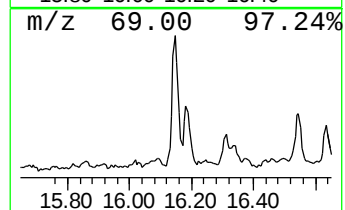
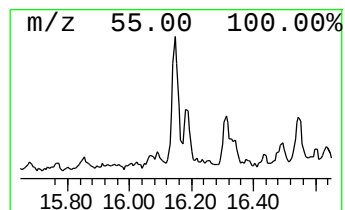
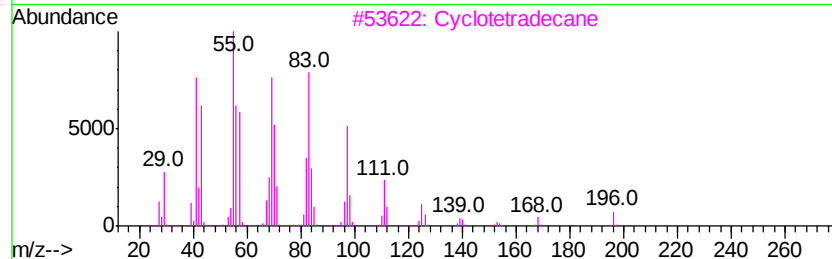
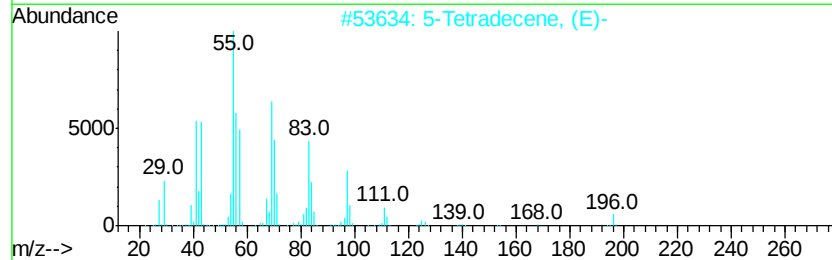
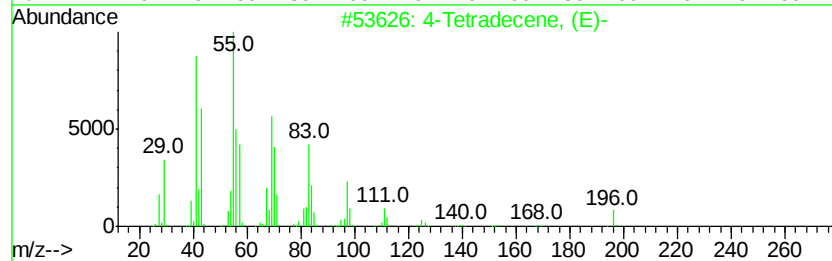
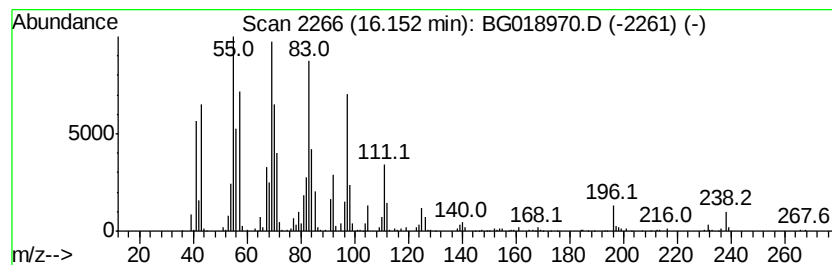
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 (DEL) Alkane: Cyclic16.15 Concentration Rank 45

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.15 | 5.51 ng/ul | 326583 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | 4-Tetradecene, (E)- | 196 | C14H28  | 041446-78-0 | 97   |
| 2    |    |   | 5-Tetradecene, (E)- | 196 | C14H28  | 041446-66-6 | 96   |
| 3    |    |   | Cyclotetradecane    | 196 | C14H28  | 000295-17-0 | 95   |
| 4    |    |   | 8-Heptadecene       | 238 | C17H34  | 054290-12-9 | 95   |
| 5    |    |   | 6-Tetradecene, (Z)- | 196 | C14H28  | 041446-61-1 | 94   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

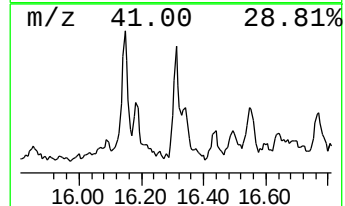
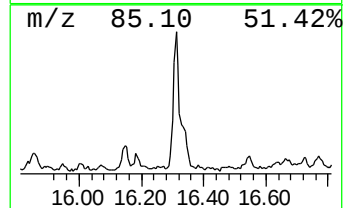
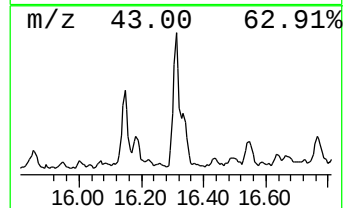
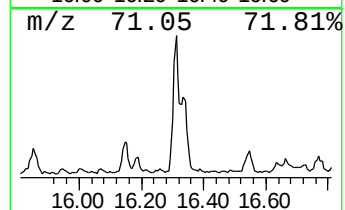
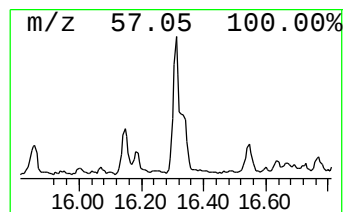
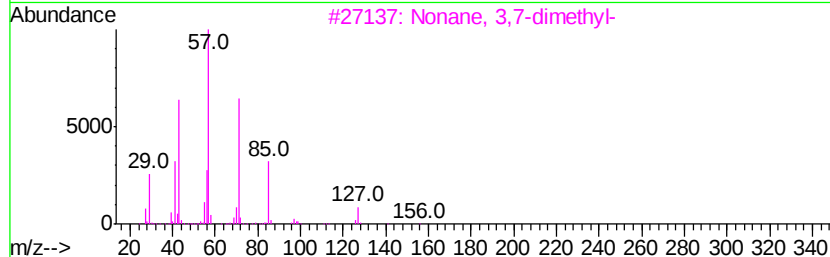
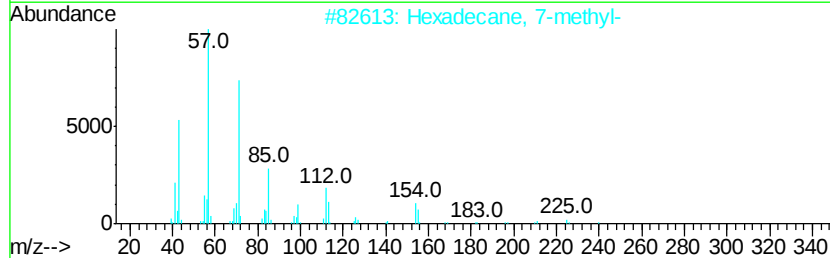
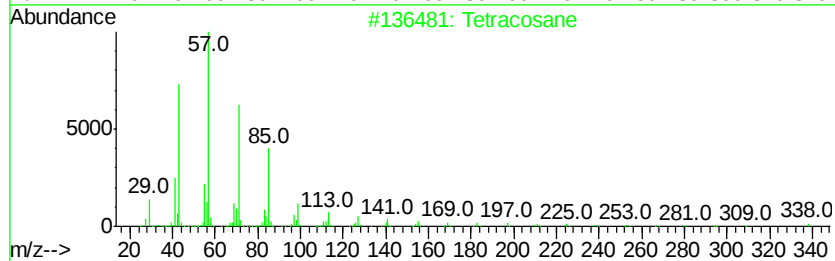
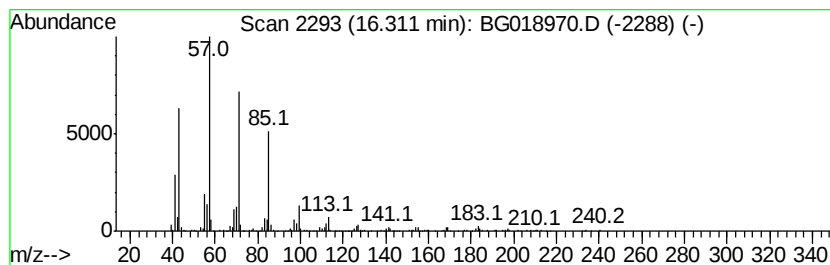
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 35

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.31 | 6.98 ng/ul | 413367 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Tetracosane           | 338 | C24H50  | 000646-31-1 | 86   |
| 2    |    |   | Hexadecane, 7-methyl- | 240 | C17H36  | 026730-20-1 | 76   |
| 3    |    |   | Nonane, 3,7-dimethyl- | 156 | C11H24  | 017302-32-8 | 74   |
| 4    |    |   | Eicosane              | 282 | C20H42  | 000112-95-8 | 74   |
| 5    |    |   | Pentadecane           | 212 | C15H32  | 000629-62-9 | 72   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

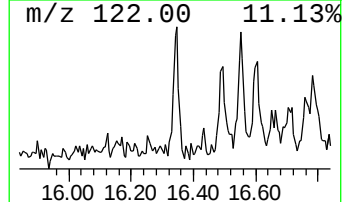
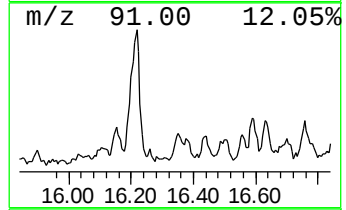
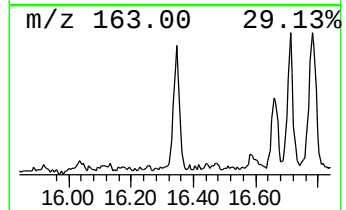
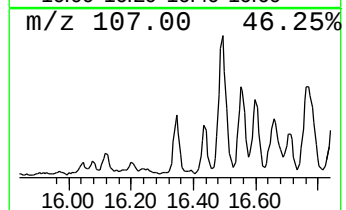
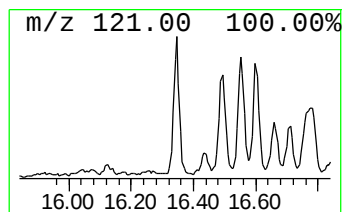
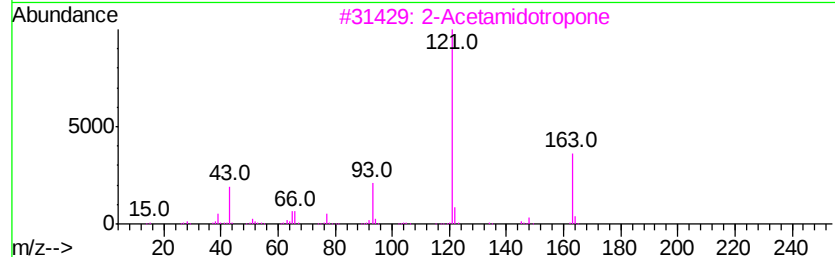
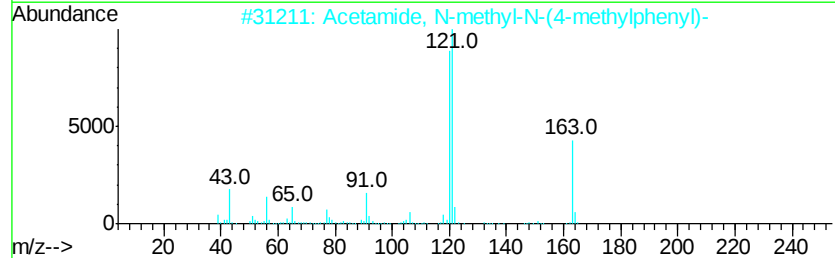
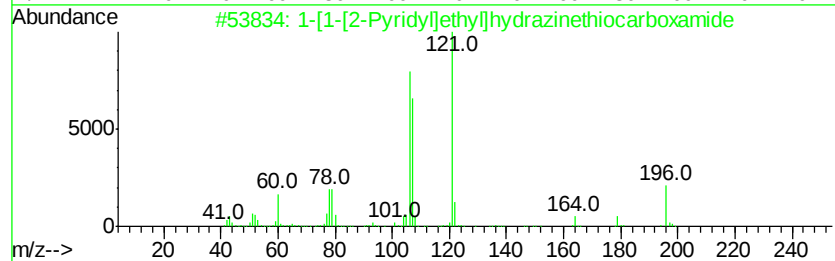
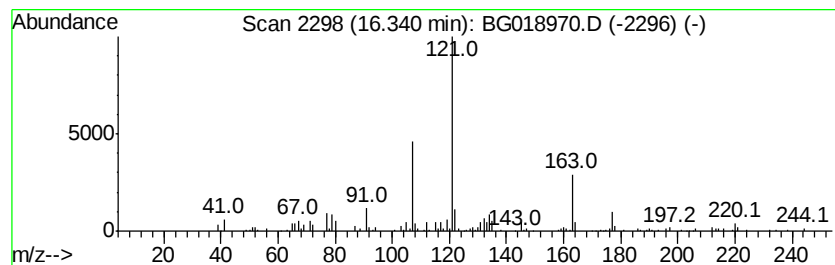
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 31 1-[1-[2-Pyridyl]ethyl]hydra... Concentration Rank 42

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.34 | 6.22 ng/ul | 368702 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1-[1-[2-Pyridyl]ethyl]hydrazinet... | 196 | C8H12N4S | 104692-01-5 | 59   |
| 2    |    |   | Acetamide, N-methyl-N-(4-methylp... | 163 | C10H13NO | 000612-03-3 | 47   |
| 3    |    |   | 2-Acetamidotropone                  | 163 | C9H9NO2  | 006422-12-4 | 47   |
| 4    |    |   | Benzene, 1-methoxy-4-octyl-         | 220 | C15H24O  | 067698-82-2 | 43   |
| 5    |    |   | p-Sec-butylphenyl acetate           | 192 | C12H16O2 | 003245-24-7 | 43   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

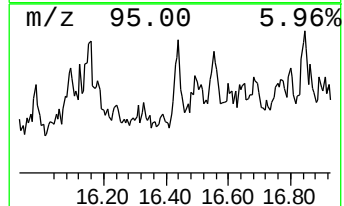
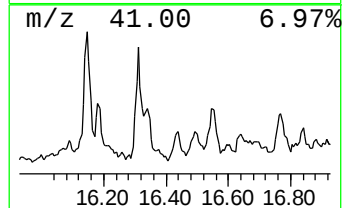
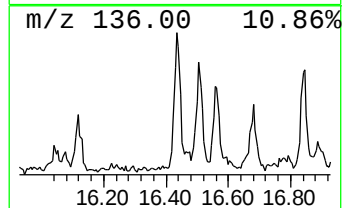
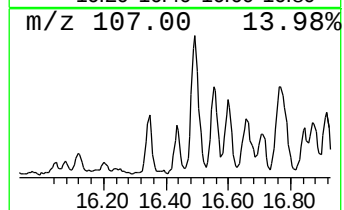
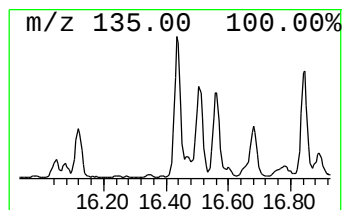
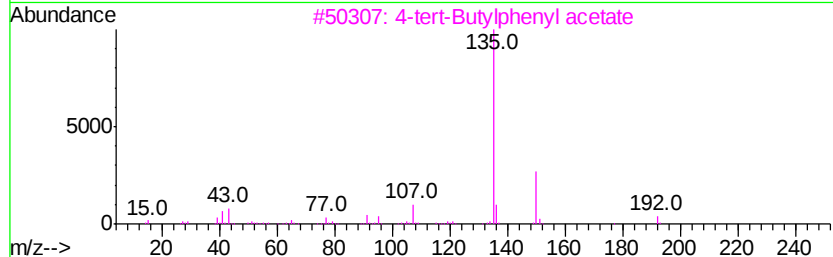
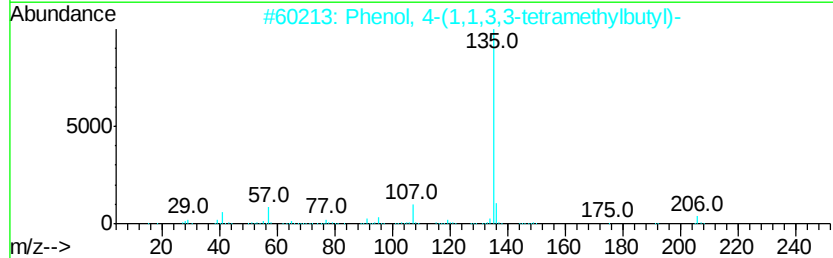
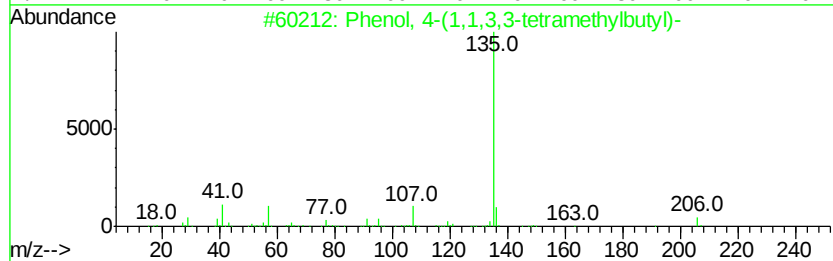
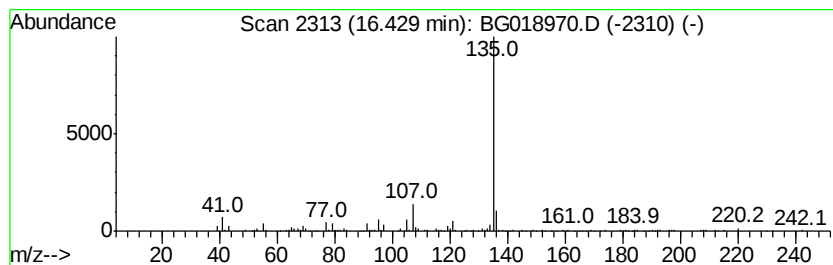
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 65

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.43 | 3.31 ng/ul | 196154 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O  | 000140-66-9 | 72   |
| 2    |    |   | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O  | 000140-66-9 | 72   |
| 3    |    |   | 4-tert-Butylphenyl acetate          | 192 | C12H16O2 | 003056-64-2 | 72   |
| 4    |    |   | Phenol, 2-methyl-5-(1-methylethyl)- | 150 | C10H14O  | 000499-75-2 | 72   |
| 5    |    |   | 2-Methyl-2-(4'-hydroxyphenyl)pen... | 192 | C12H16O2 | 070205-18-4 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

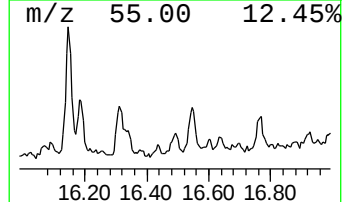
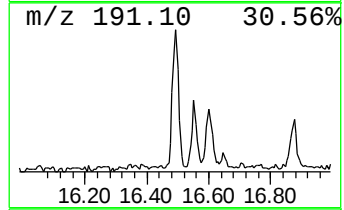
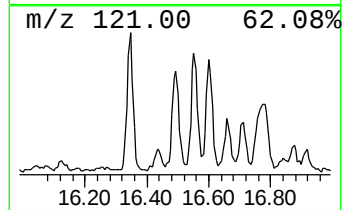
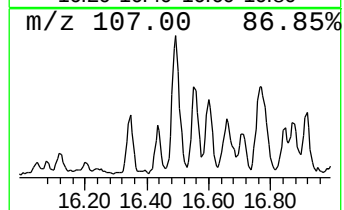
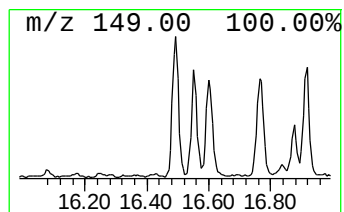
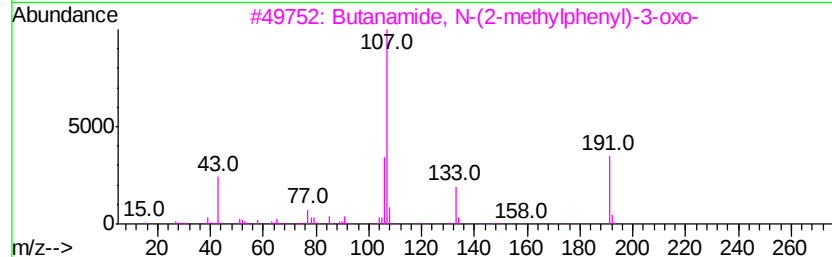
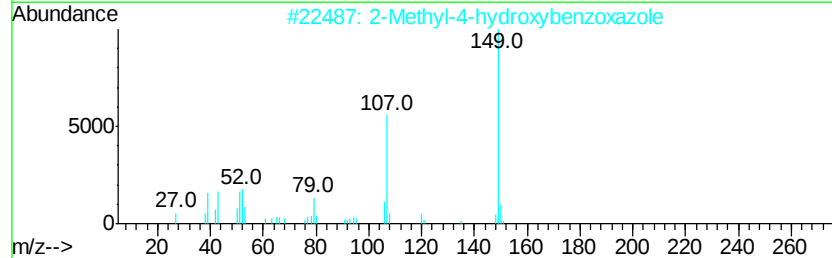
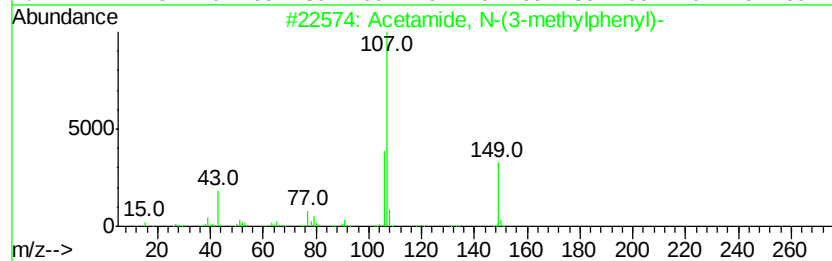
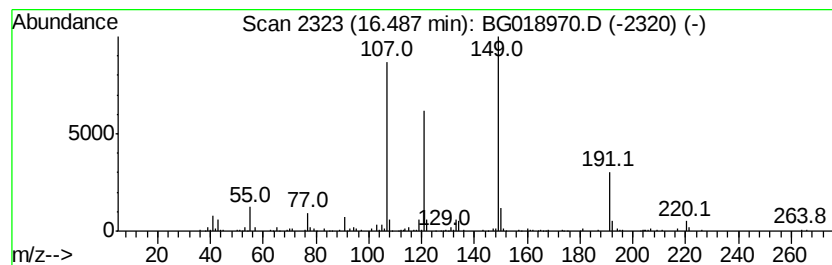
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 Acetamide, N-(3-methylphenyl)- Concentration Rank 58

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.49 | 3.65 ng/ul | 216255 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID                                                | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Acetamide, N-(3-methylphenyl)-                              | 149 | C9H11NO   | 000537-92-8 | 53   |
| 2    |    |   | 2-Methyl-4-hydroxybenzoxazole                               | 149 | C8H7NO2   | 051110-60-2 | 53   |
| 3    |    |   | Butanamide, N-(2-methylphenyl)-3-oxo-                       | 191 | C11H13NO2 | 000093-68-5 | 38   |
| 4    |    |   | Pyridine-3-carbonitrile, 1,2-dihydro-                       | 191 | C9H9N3O2  | 220098-92-0 | 38   |
| 5    |    |   | Phenol, 2-methyl-4-(1,1,3,3-tetrahydro-1H-benzoxazol-2-yl)- | 220 | C15H24O   | 002219-84-3 | 30   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

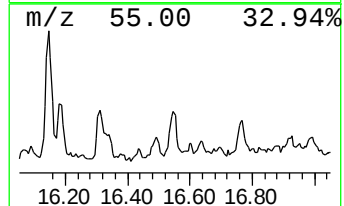
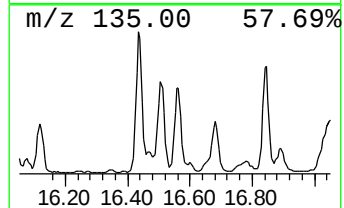
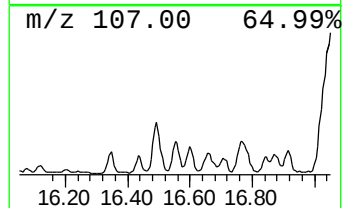
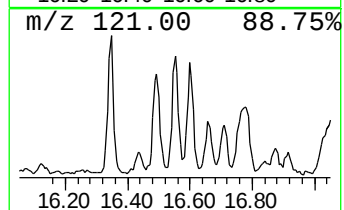
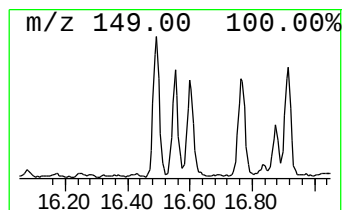
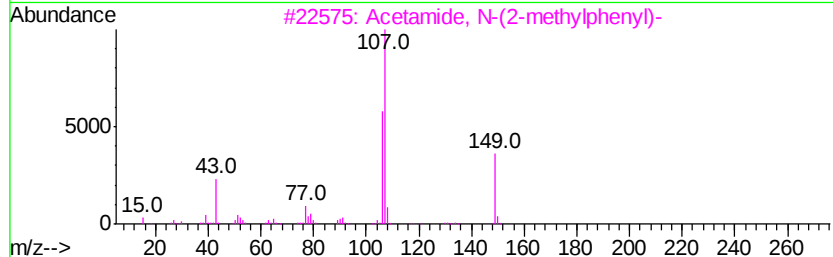
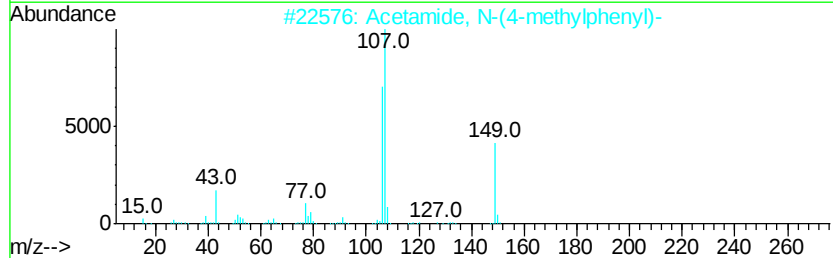
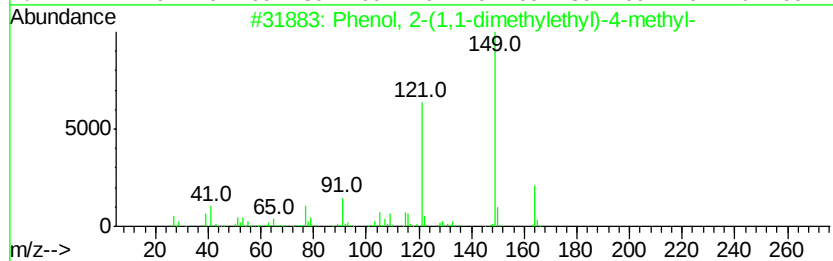
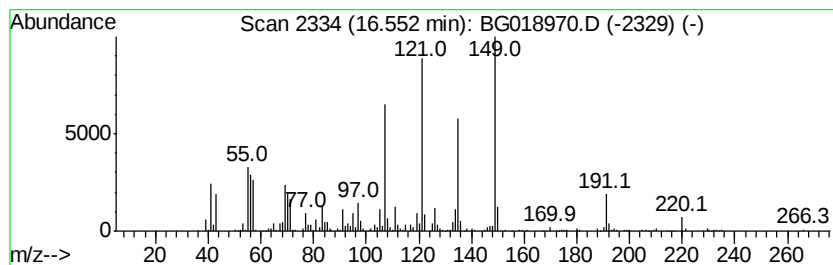
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 unknown-03 Concentration Rank 52

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.55 | 4.87 ng/ul | 288350 | Phenanthrene-d10 | 17.37 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | Phenol, 2-(1,1-dimethylethyl)-4-... | 164 | C11H16O      | 002409-55-4 | 43      |      |      |
| 2    | Acetamide, N-(4-methylphenyl)-      | 149 | C9H11NO      | 000103-89-9 | 30      |      |      |
| 3    | Acetamide, N-(2-methylphenyl)-      | 149 | C9H11NO      | 000120-66-1 | 30      |      |      |
| 4    | Acetamide, N-(2-methylphenyl)-      | 149 | C9H11NO      | 000120-66-1 | 27      |      |      |
| 5    | Phenol, 2-methyl-4-(1,1,3,3-tetr... | 220 | C15H24O      | 002219-84-3 | 25      |      |      |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

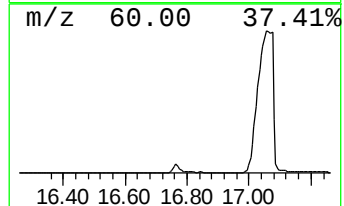
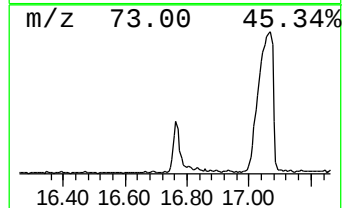
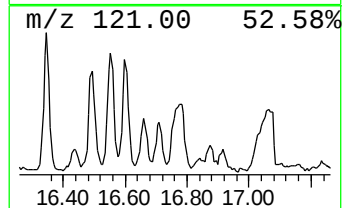
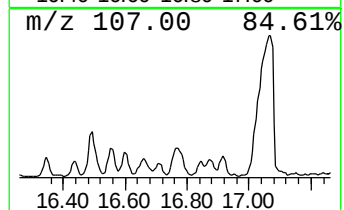
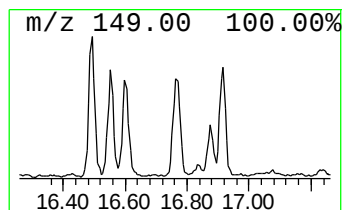
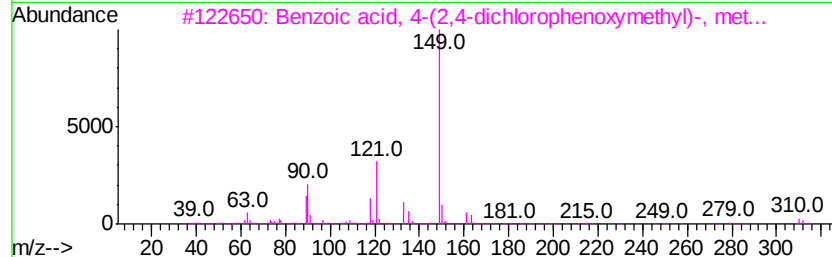
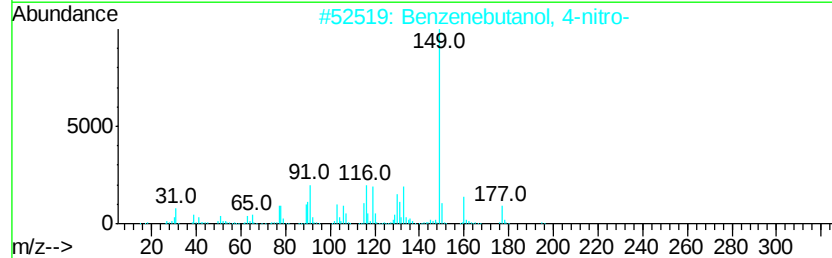
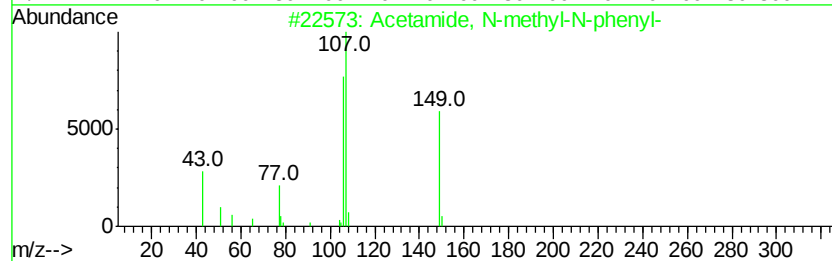
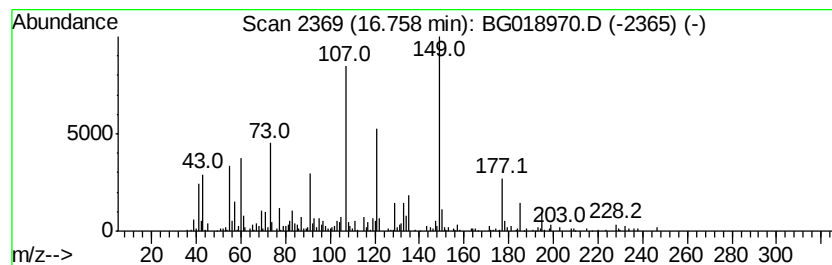
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 unknown-04 Concentration Rank 44

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.76 | 5.74 ng/ul | 340418 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Acetamide, N-methyl-N-phenyl-       | 149 | C9H11NO     | 000579-10-2  | 38   |
| 2    |    |   | Benzenebutanol, 4-nitro-            | 195 | C10H13NO3   | 079524-20-2  | 35   |
| 3    |    |   | Benzoic acid, 4-(2,4-dichlorophe... | 310 | C15H12Cl2O3 | 1000294-38-1 | 35   |
| 4    |    |   | 1,8-Nonadiene, 2,7-dimethyl-5-(1... | 192 | C14H24      | 068702-20-5  | 30   |
| 5    |    |   | Rose acetate                        | 266 | C10H9Cl3O2  | 1000109-21-6 | 30   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

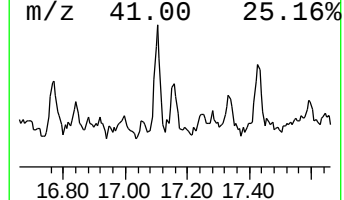
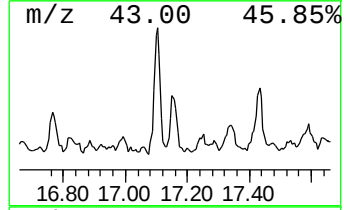
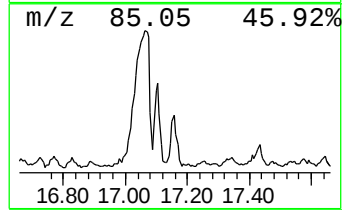
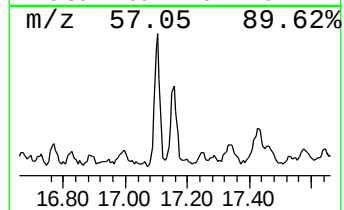
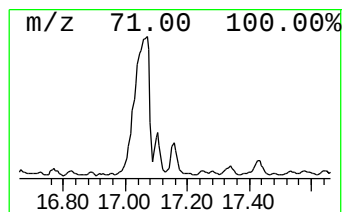
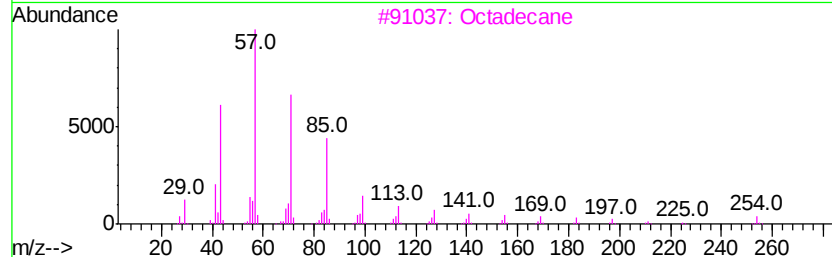
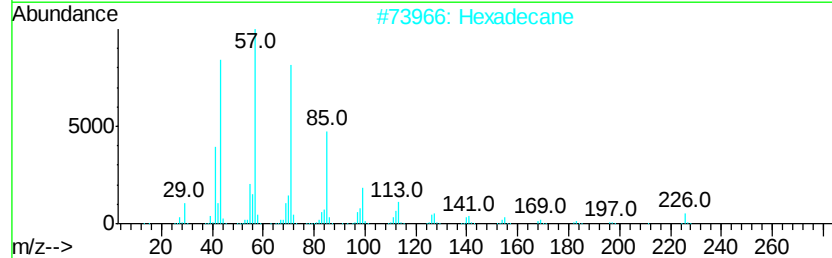
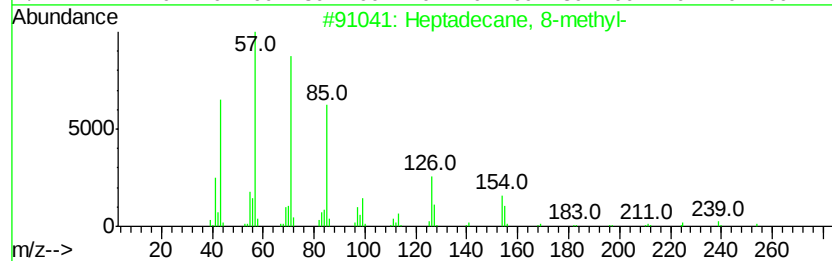
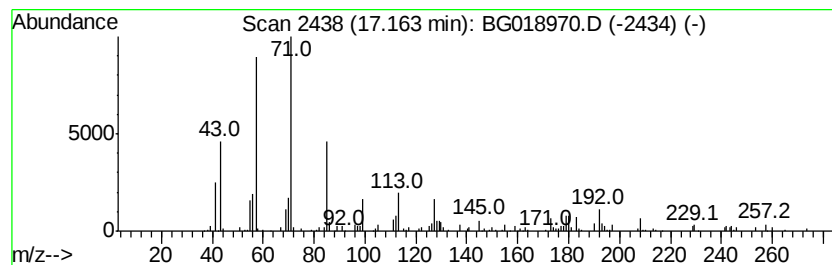
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 54

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.16 | 4.47 ng/ul | 265043 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptadecane, 8-methyl- | 254 | C18H38  | 013287-23-5 | 86   |
| 2    |    | Hexadecane             | 226 | C16H34  | 000544-76-3 | 86   |
| 3    |    | Octadecane             | 254 | C18H38  | 000593-45-3 | 86   |
| 4    |    | Hexadecane             | 226 | C16H34  | 000544-76-3 | 78   |
| 5    |    | Tetracosane            | 338 | C24H50  | 000646-31-1 | 78   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

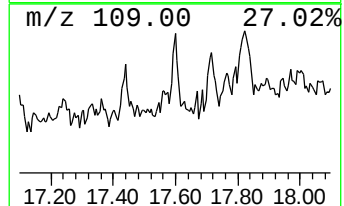
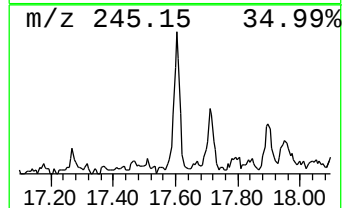
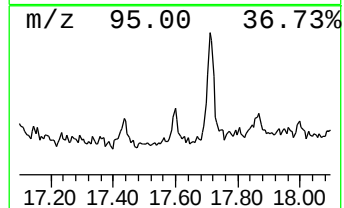
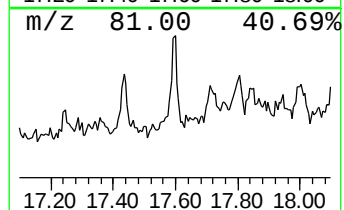
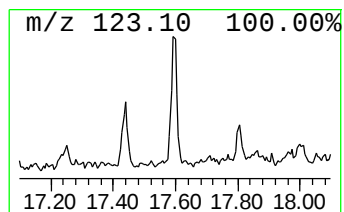
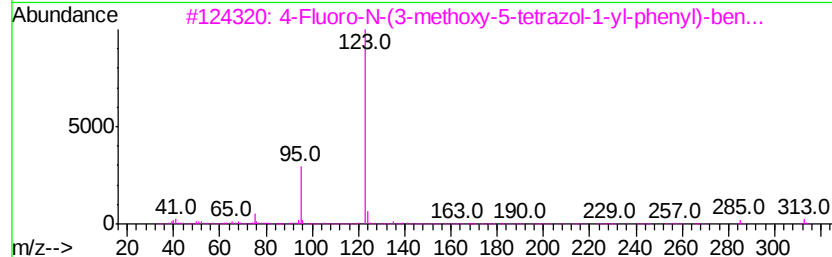
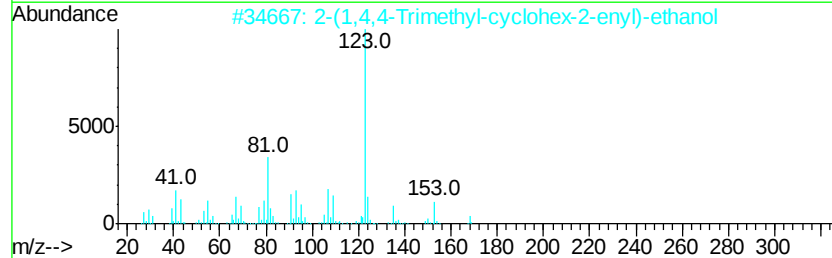
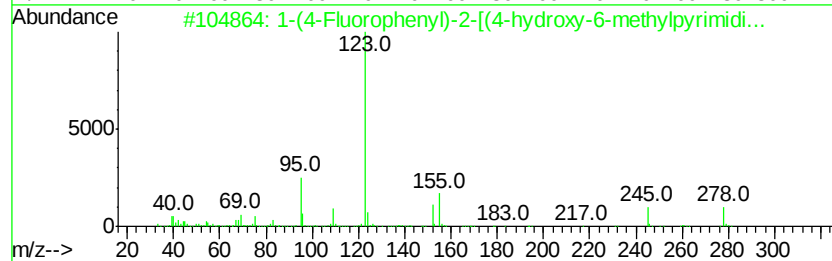
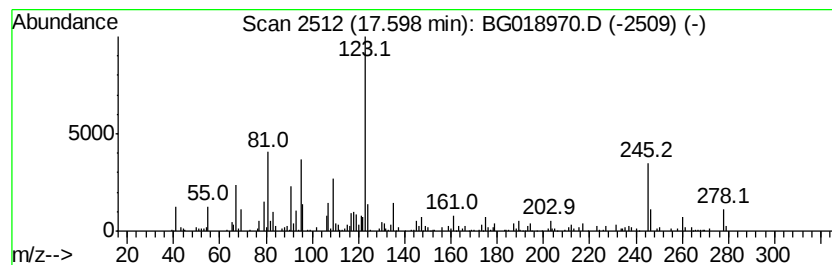
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 37 1-(4-Fluorophenyl)-2-[(4-hydroxy... Concentration Rank 74

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.60 | 2.27 ng/ul | 134570 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID                         | MW  | MolForm      | CAS#         | Qual |
|------|----|--------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | 1-(4-Fluorophenyl)-2-[(4-hydroxy...  | 278 | C13H11FN2O2S | 119730-11-9  | 62   |
| 2    |    | 2-(1,4,4-Trimethyl-cyclohex-2-en...  | 168 | C11H20O      | 1000190-92-3 | 47   |
| 3    |    | 4-Fluoro-N-(3-methoxy-5-tetrazol...  | 313 | C15H12FN5O2  | 1000295-31-2 | 47   |
| 4    |    | 1-(1-Ethyl-2,3-dimethyl-cyclopent... | 166 | C11H18O      | 1000185-18-6 | 47   |
| 5    |    | Cyclopropanecarboxylic acid, 2,2...  | 328 | C21H28O3     | 000121-21-1  | 43   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

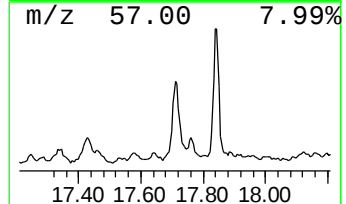
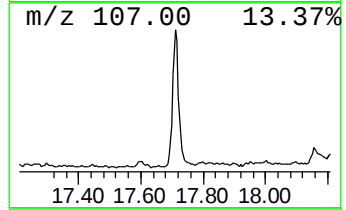
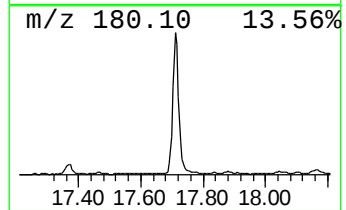
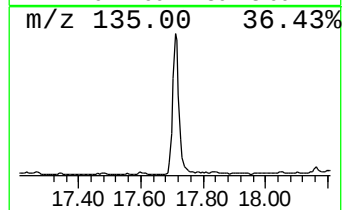
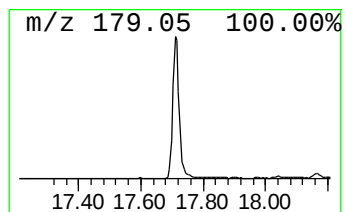
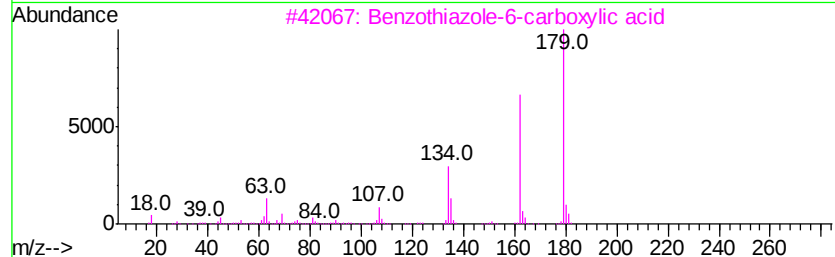
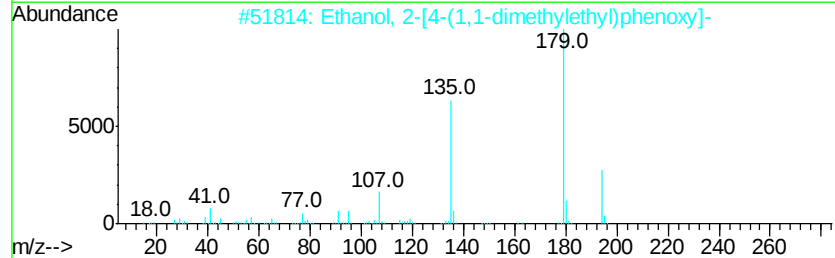
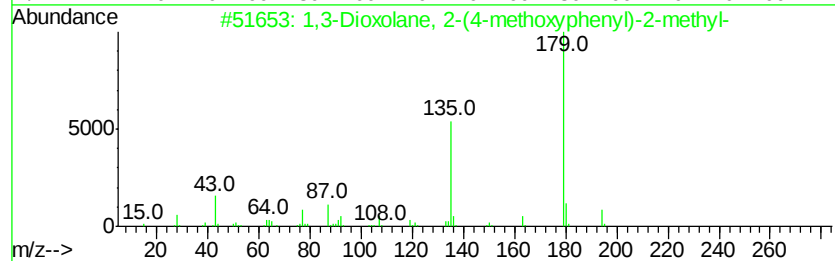
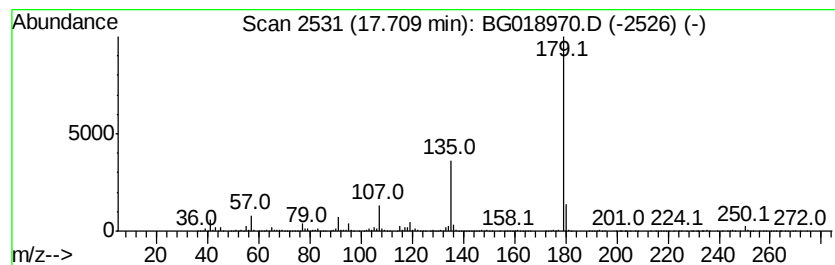
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 38 1,3-Dioxolane, 2-(4-methoxy... Concentration Rank 5

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.71 | 26.22 ng/ul | 1553990 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID                                 | MW  | MolForm   | CAS#        | Qual |
|------|----|----------------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 1,3-Dioxolane, 2-(4-methoxyphenyl)-2-methyl- | 194 | C11H14O3  | 036881-00-2 | 56   |
| 2    |    | Ethanol, 2-[4-(1,1-dimethylethyl)phenoxy]-   | 194 | C12H18O2  | 000713-46-2 | 50   |
| 3    |    | Benzothiazole-6-carboxylic acid              | 179 | C8H5N02S  | 003622-35-3 | 50   |
| 4    |    | Ethanol, 2-[4-(1,1-dimethylpropyl)phenoxy]-  | 208 | C13H20O2  | 006382-07-6 | 43   |
| 5    |    | 1-Dimethyl(phenyl)silyloxypropane            | 194 | C11H18OSi | 013360-21-9 | 39   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

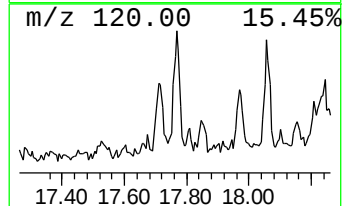
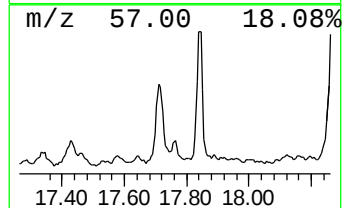
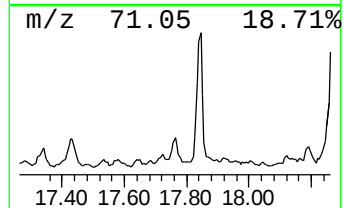
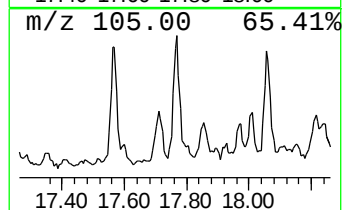
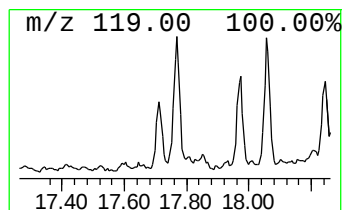
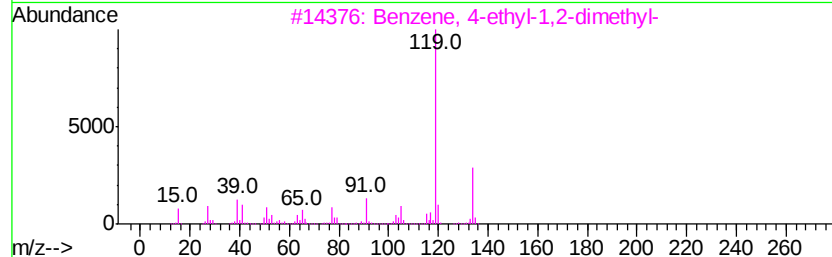
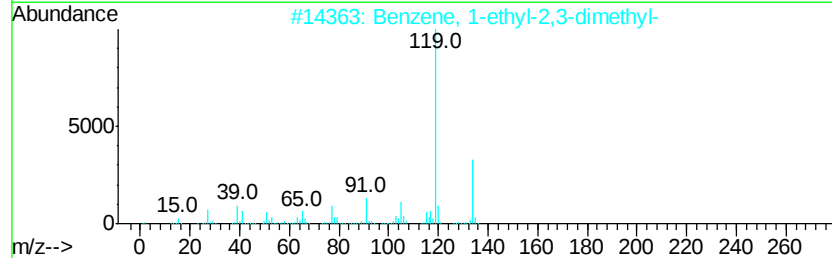
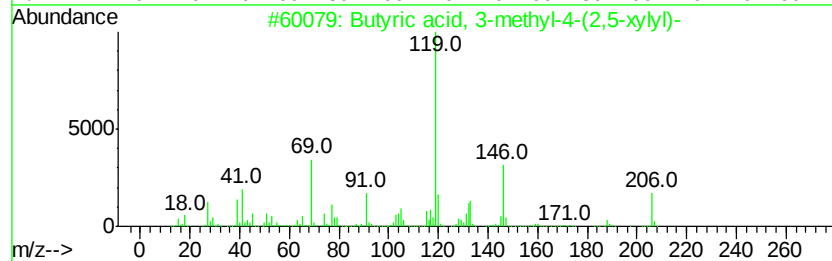
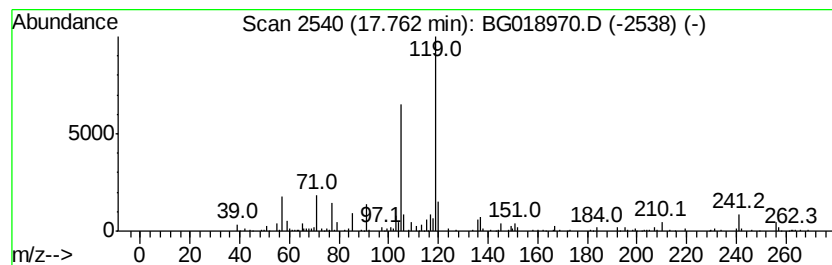
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 Butyric acid, 3-methyl-4-(2... Concentration Rank 67

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.76 | 3.06 ng/ul | 181591 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Butyric acid, 3-methyl-4-(2,5-xv... | 206 | C13H18O2 | 030275-76-4 | 59   |
| 2    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14   | 000933-98-2 | 58   |
| 3    |    | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14   | 000934-80-5 | 53   |
| 4    |    | Benzene, 1-methyl-4-(1-methyleth... | 134 | C10H14   | 000099-87-6 | 53   |
| 5    |    | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14   | 000934-74-7 | 53   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

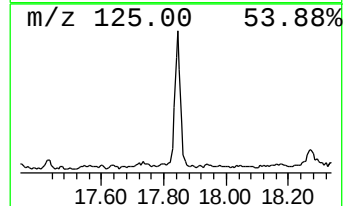
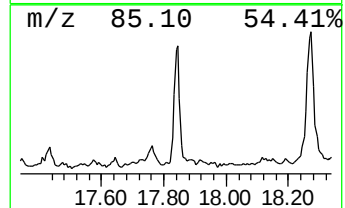
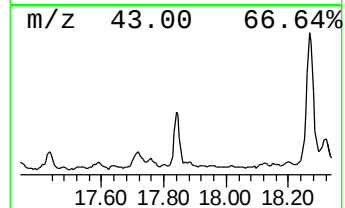
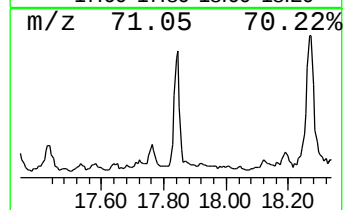
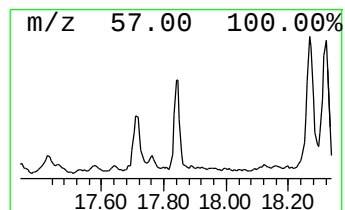
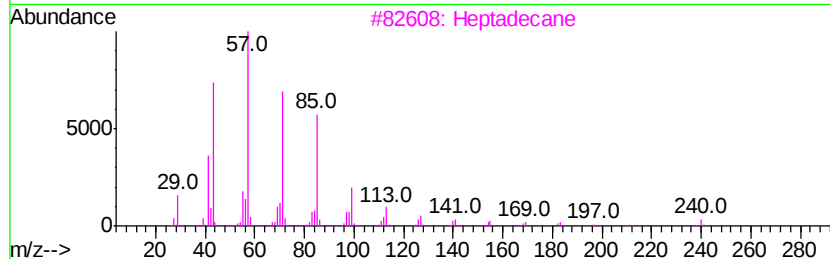
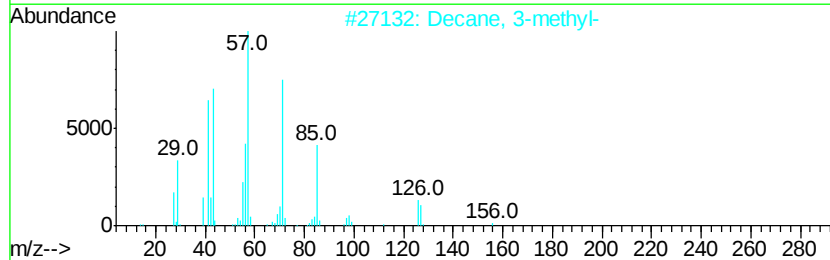
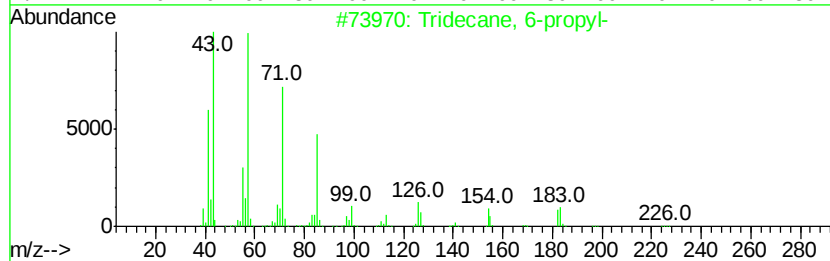
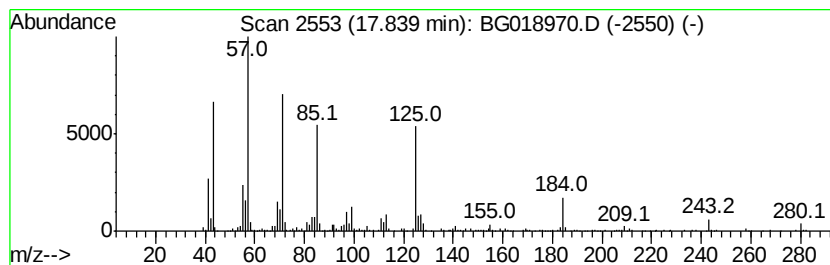
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 17.84 | 12.12 ng/ul | 718348 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane, 6-propyl- | 226 | C16H34  | 055045-10-8 | 78   |
| 2    |    | Decane, 3-methyl-    | 156 | C11H24  | 013151-34-3 | 55   |
| 3    |    | Heptadecane          | 240 | C17H36  | 000629-78-7 | 49   |
| 4    |    | Heneicosane          | 296 | C21H44  | 000629-94-7 | 49   |
| 5    |    | Undecane, 2-methyl-  | 170 | C12H26  | 007045-71-8 | 49   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

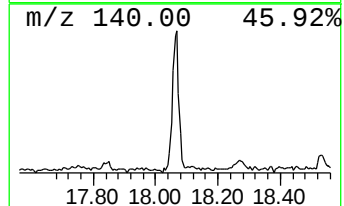
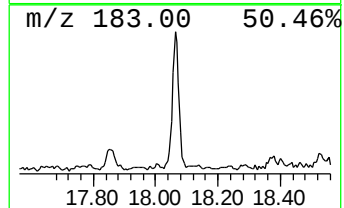
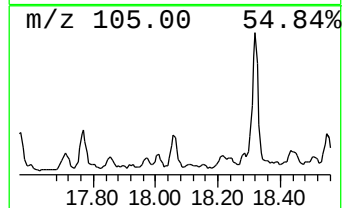
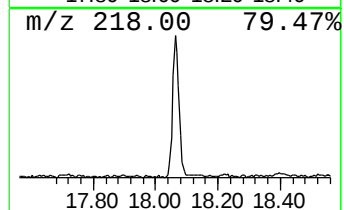
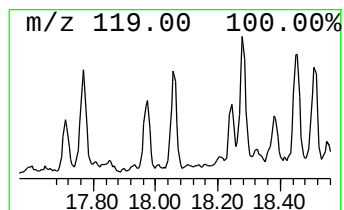
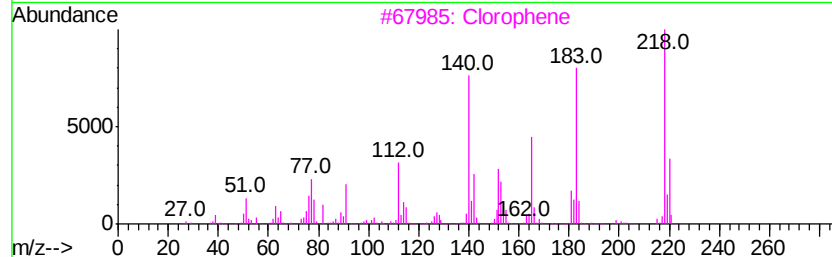
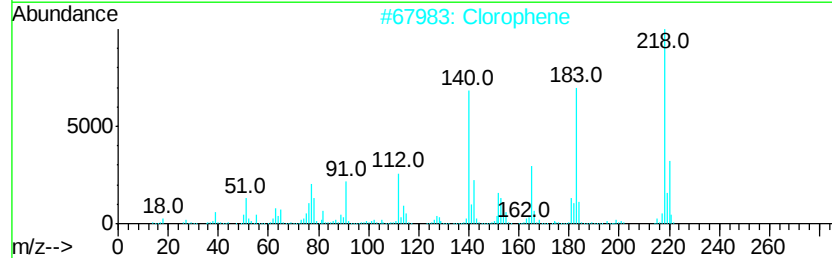
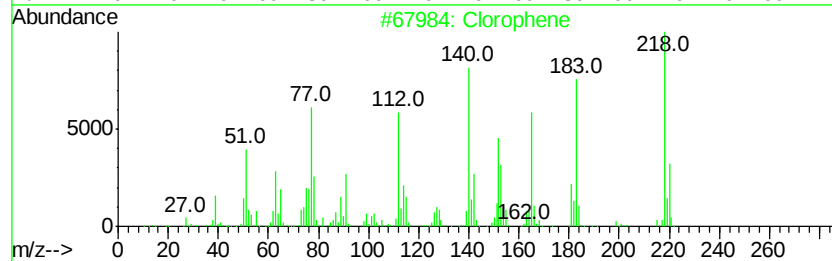
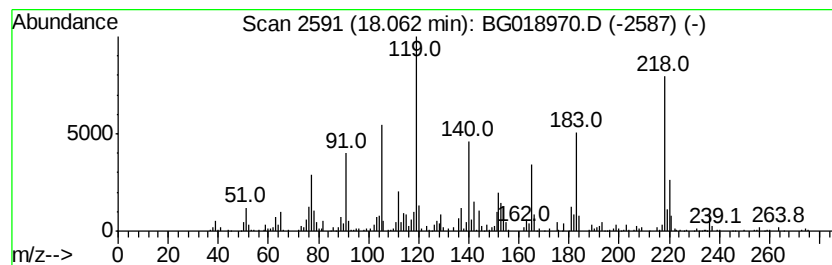
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 41 Clorophene Concentration Rank 43

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.06 | 6.07 ng/ul | 359732 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Clorophene                          | 218 | C13H11ClO  | 000120-32-1  | 99   |
| 2    |    | Clorophene                          | 218 | C13H11ClO  | 000120-32-1  | 98   |
| 3    |    | Clorophene                          | 218 | C13H11ClO  | 000120-32-1  | 98   |
| 4    |    | Clorophene acetate                  | 260 | C15H13ClO2 | 1000120-34-0 | 58   |
| 5    |    | Benzenesulfonyl chloride, 2,4,6-... | 218 | C9H11ClO2S | 000773-64-8  | 52   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

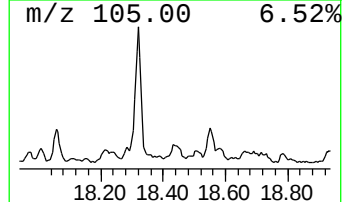
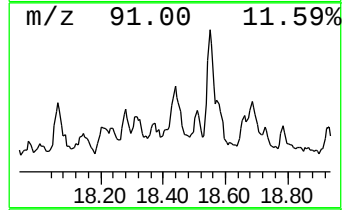
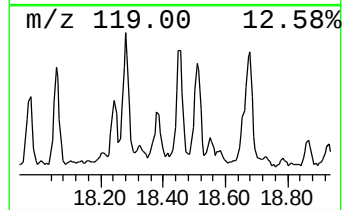
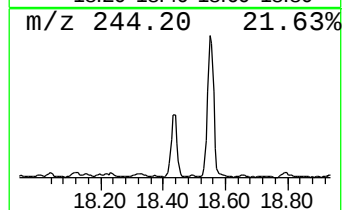
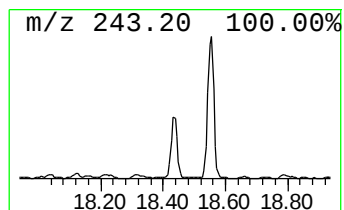
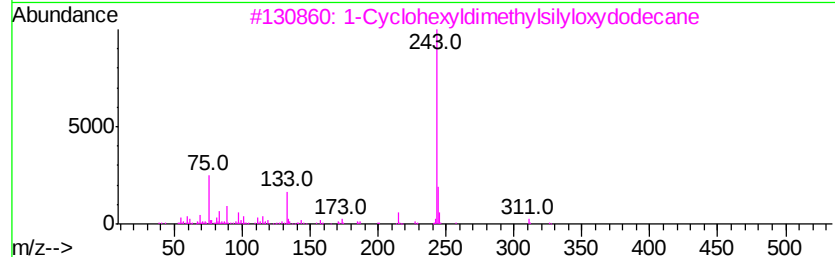
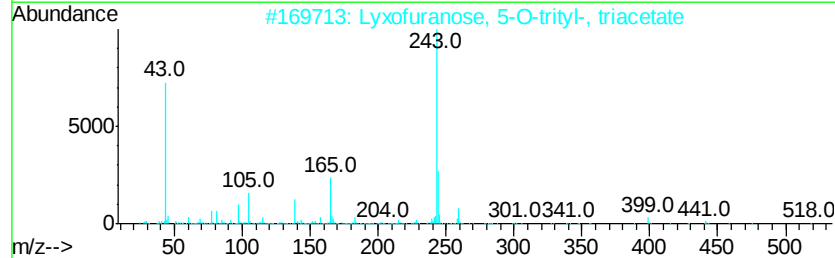
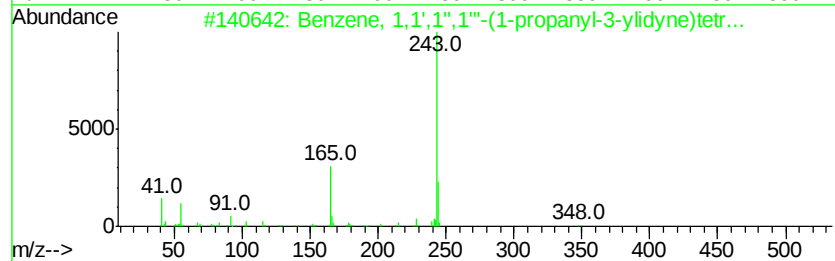
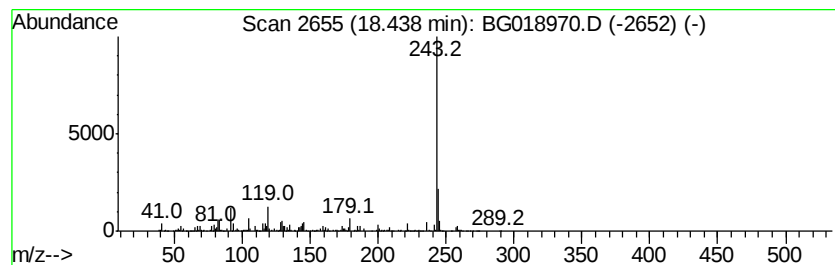
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 42 Benzene, 1,1',1'',1'''-(1-p... Concentration Rank 25

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 18.44 | 10.49 ng/ul | 621461 | Phenanthrene-d10 | 17.37 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|---------|---|-------------------------------------|-----|-----------|--------------|------|
| 1       |   | Benzene, 1,1',1'',1'''-(1-propan... | 348 | C27H24    | 054966-03-9  | 53   |
| 2       |   | Lyxofuranose, 5-O-trityl-, triac... | 518 | C30H30O8  | 103419-92-7  | 45   |
| 3       |   | 1-Cyclohexyldimethylsilyloxvode...  | 326 | C20H42OSi | 1000281-96-3 | 45   |
| 4       |   | 3-Phenyl-4-azafluorene              | 243 | C18H13N   | 033777-97-8  | 42   |
| 5       |   | 1-Tritylaziridine-2-carbothioic ... | 421 | C28H23NOS | 1000191-60-9 | 42   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

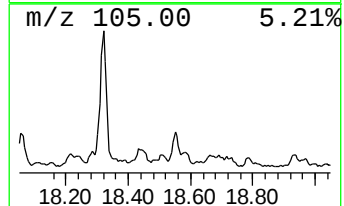
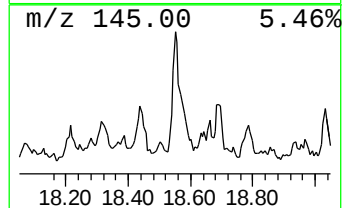
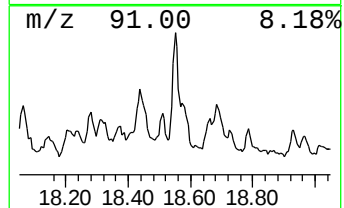
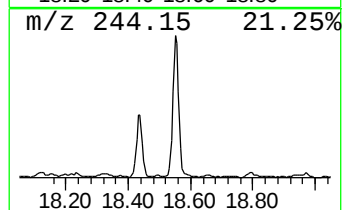
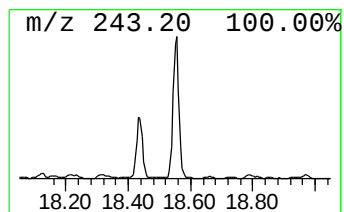
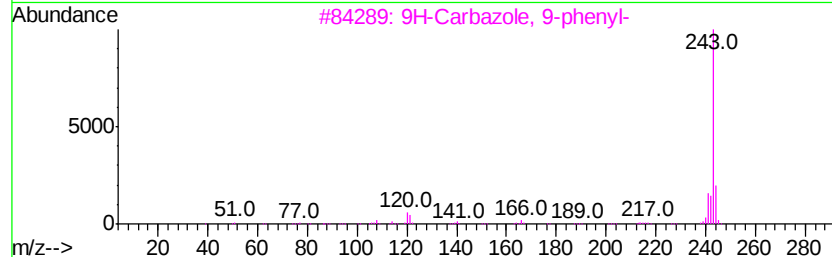
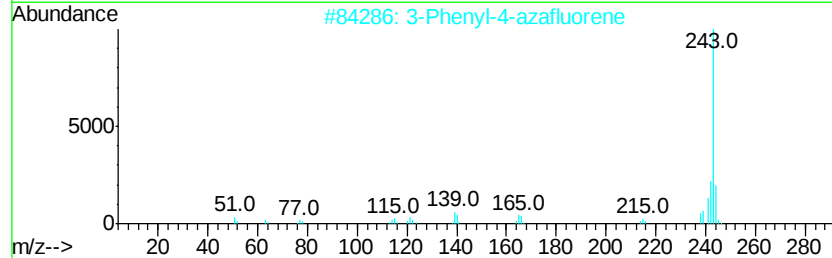
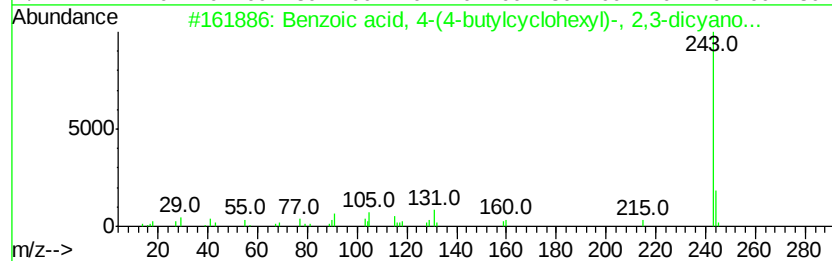
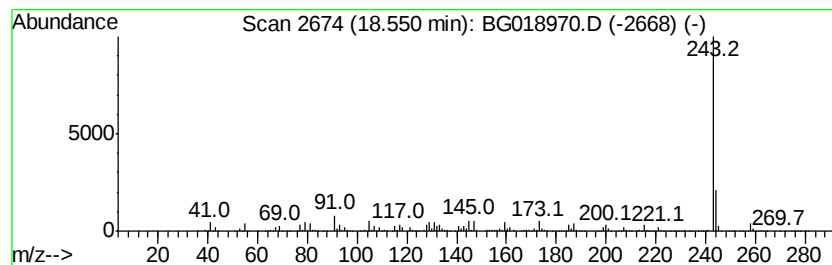
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 43 Benzoic acid, 4-(4-butylcyclohexyl) Concentration Rank 12

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.55 | 17.86 ng/ul | 1058170 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Benzoic acid, 4-(4-butylcyclohexyl) | 430 | C27H30N2O3 | 086377-40-4 | 64   |
| 2    |    | 3-Phenyl-4-azafluorene              | 243 | C18H13N    | 033777-97-8 | 59   |
| 3    |    | 9H-Carbazole, 9-phenyl-             | 243 | C18H13N    | 001150-62-5 | 59   |
| 4    |    | 9H-Carbazole, 9-phenyl-             | 243 | C18H13N    | 001150-62-5 | 59   |
| 5    |    | 1,1,1-triphenyl-2-decanone          | 384 | C28H32O    | 072152-27-3 | 45   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

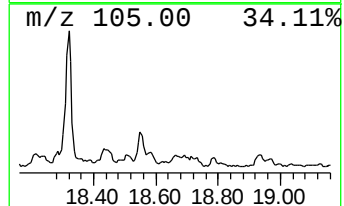
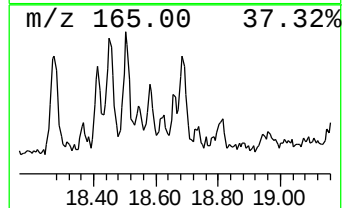
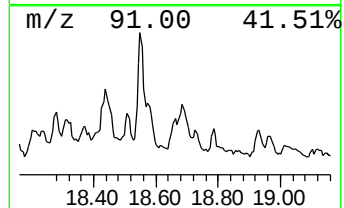
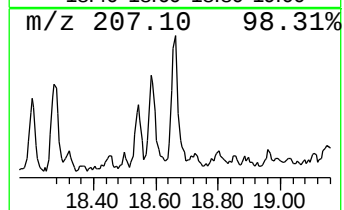
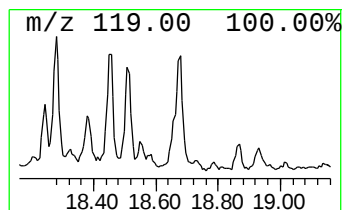
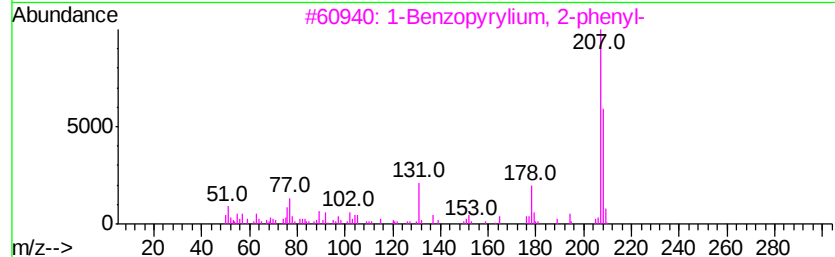
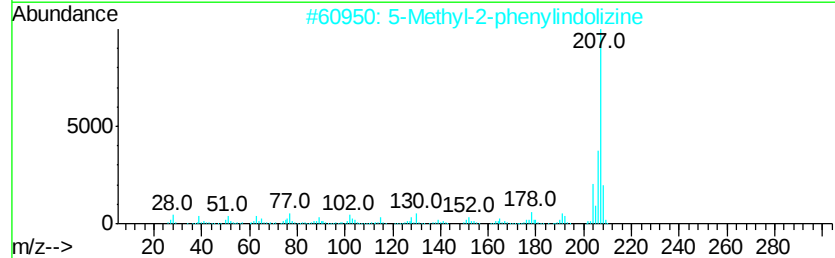
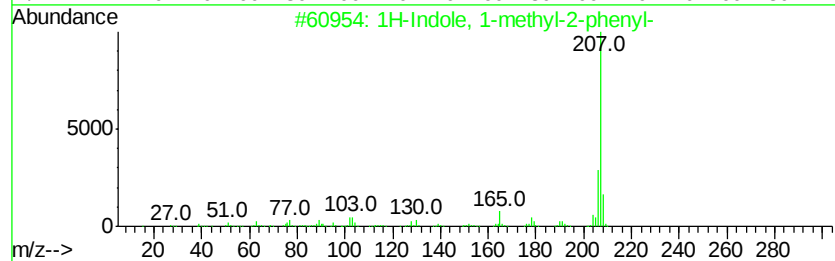
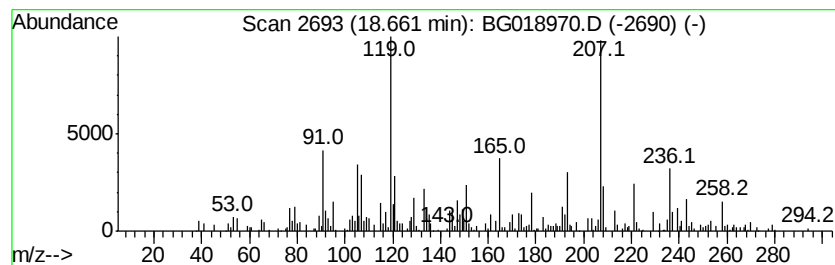
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 44 unknown-05 Concentration Rank 73

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.66 | 2.31 ng/ul | 136748 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1H-Indole, 1-methyl-2-phenyl-        | 207 | C15H13N  | 003558-24-5 | 25   |
| 2    |    |   | 5-Methyl-2-phenylindolizine          | 207 | C15H13N  | 036944-99-7 | 25   |
| 3    |    |   | 1-Benzopyrvium, 2-phenyl-            | 207 | C15H11O  | 014051-53-7 | 25   |
| 4    |    |   | 1H-Indole, 5-methyl-2-phenyl-        | 207 | C15H13N  | 013228-36-9 | 18   |
| 5    |    |   | Benzenepropanoic acid, .beta., .b... | 178 | C11H14O2 | 001010-48-6 | 18   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

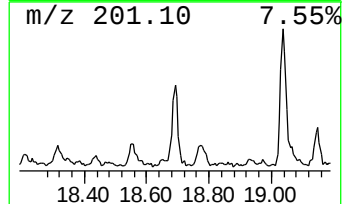
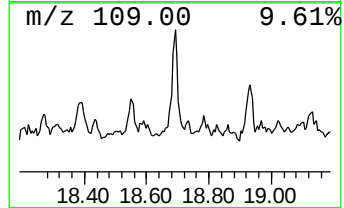
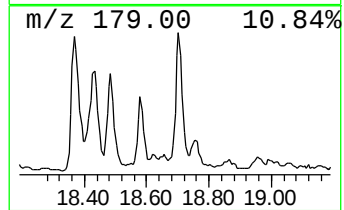
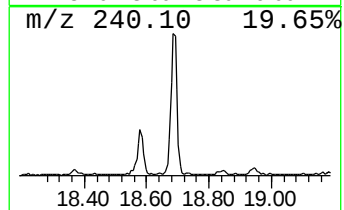
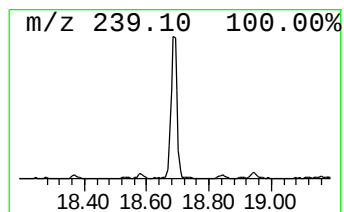
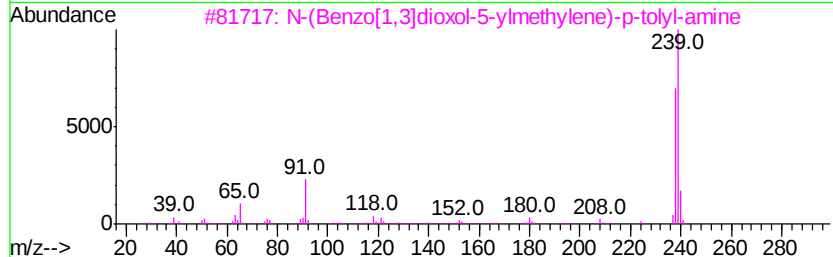
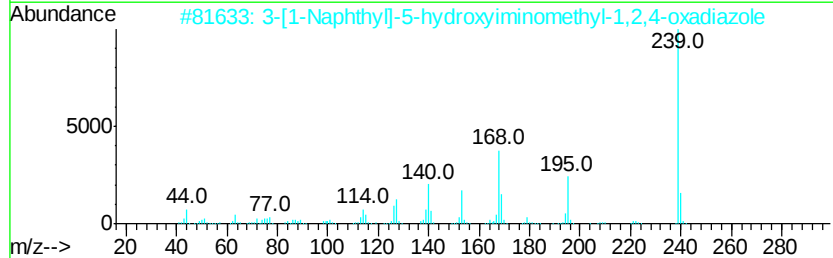
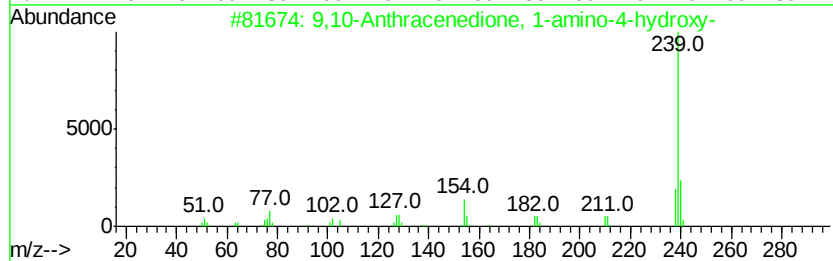
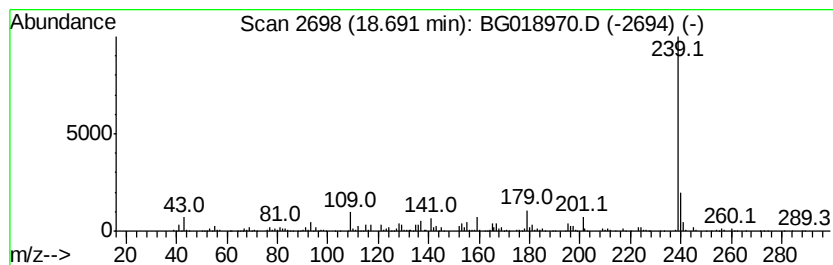
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 45 9,10-Anthracenedione, 1-ami... Concentration Rank 18

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 18.69 | 12.74 ng/ul | 755061 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID                                   | MW  | MolForm   | CAS#        | Qual |
|------|----|------------------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 9,10-Anthracenedione, 1-amino-4-...            | 239 | C14H9NO3  | 000116-85-8 | 59   |
| 2    |    | 3-[1-Naphthyl]-5-hydroxymethyl-2,4-oxadiazole  | 239 | C13H9N3O2 | 103499-04-3 | 59   |
| 3    |    | N-(Benzo[1,3]dioxol-5-yl)methyl-2,4-oxadiazole | 239 | C15H13NO2 | 007402-58-6 | 45   |
| 4    |    | Indole, 3-imidazo[2,1-b]thiazole               | 239 | C13H9N3S  | 292855-05-1 | 45   |
| 5    |    | Imidazo[2,1-b]thiazole, 5-(3-i...              | 239 | C13H9N3S  | 327097-37-0 | 42   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

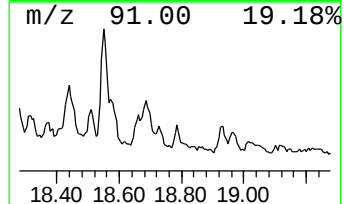
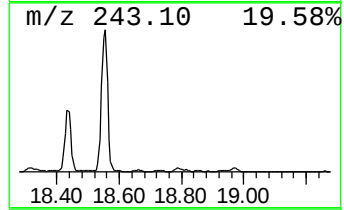
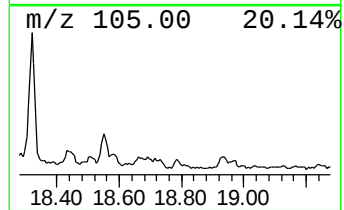
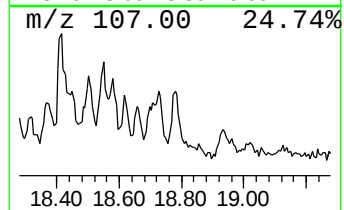
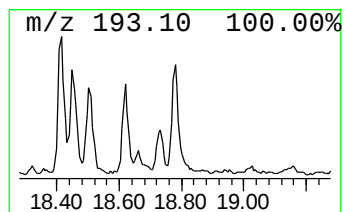
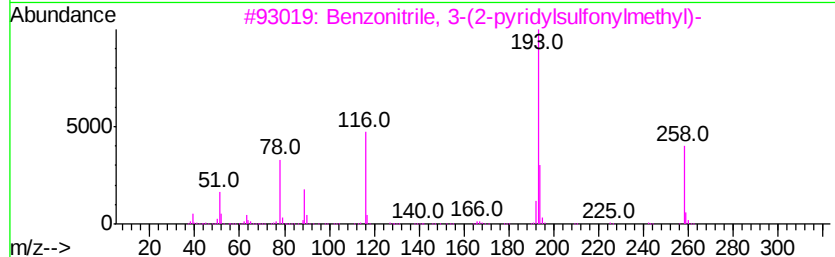
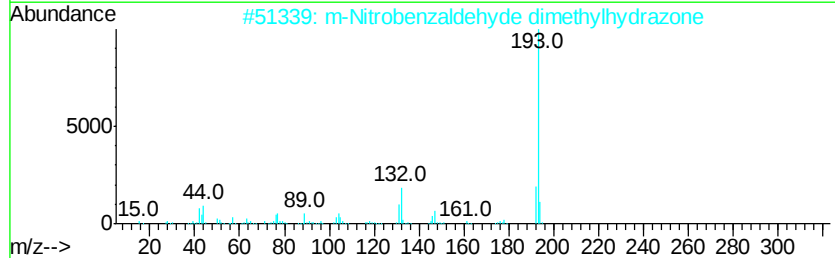
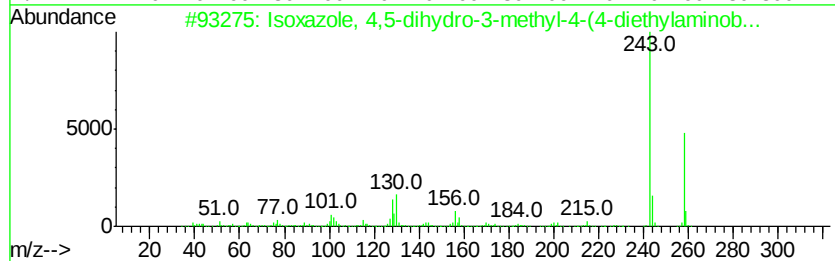
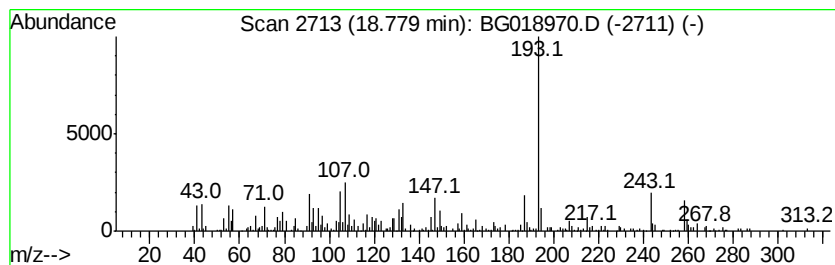
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 46 unknown-06 Concentration Rank 68

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.78 | 3.06 ng/ul | 181414 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Isoxazole, 4,5-dihydro-3-methyl-... | 258 | C15H18N2O2  | 328002-66-0  | 25   |
| 2    |    |   | m-Nitrobenzaldehyde dimethylhydr... | 193 | C9H11N3O2   | 032787-76-1  | 22   |
| 3    |    |   | Benzonitrile, 3-(2-pyridylsulfon... | 258 | C13H10N2O2S | 1000269-35-2 | 22   |
| 4    |    |   | Pyrazole, 5-methyl-3-(5-nitro-2-... | 193 | C8H7N3O3    | 016239-90-0  | 22   |
| 5    |    |   | Pentamethylnitrobenzene             | 193 | C11H15N02   | 013171-59-0  | 22   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

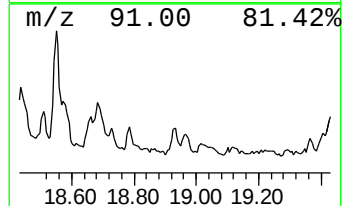
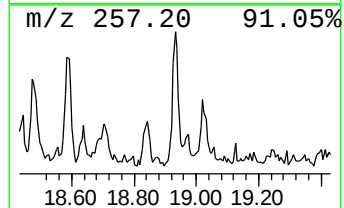
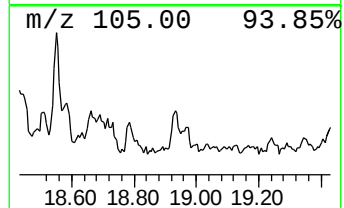
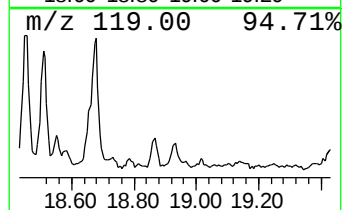
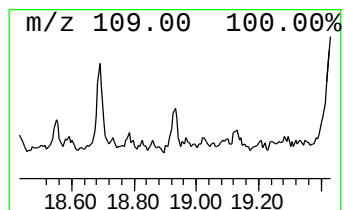
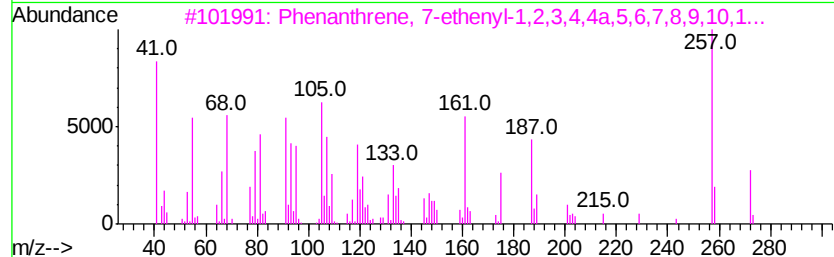
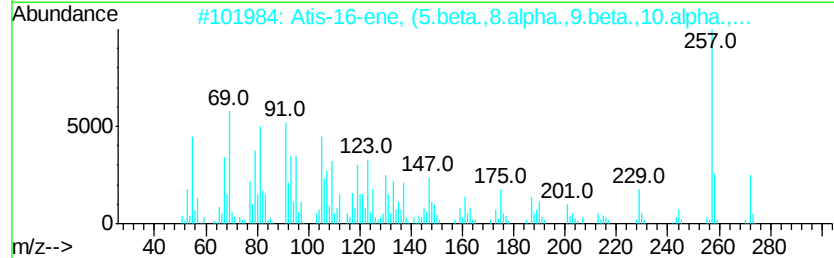
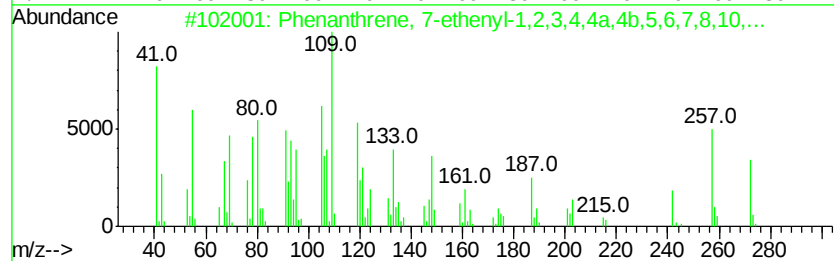
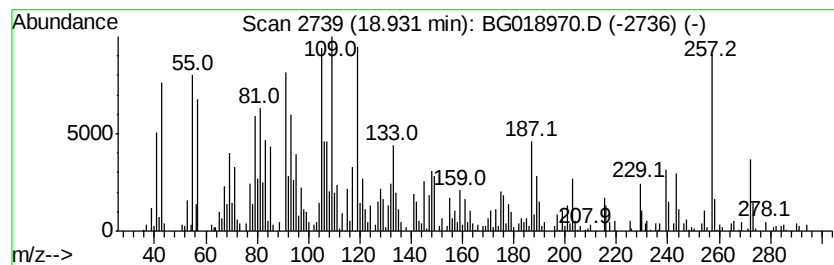
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 47 Phenanthrene, 7-ethenyl-1,2... Concentration Rank 59

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.93 | 3.58 ng/ul | 212388 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 7-ethenyl-1,2,3,4,... | 272 | C20H32  | 001686-66-4 | 58   |
| 2    |    |   | Atis-16-ene, (5.beta.,8.alpha.,9... | 272 | C20H32  | 020230-48-2 | 51   |
| 3    |    |   | Phenanthrene, 7-ethenyl-1,2,3,4,... | 272 | C20H32  | 055255-56-6 | 41   |
| 4    |    |   | Trachylobane                        | 272 | C20H32  | 005282-35-9 | 20   |
| 5    |    |   | Naphthalene, decahydro-1,1,4a-tr... | 272 | C20H32  | 000511-02-4 | 15   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

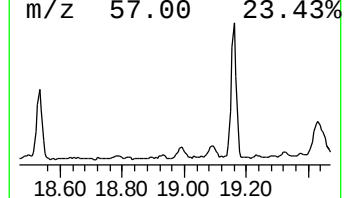
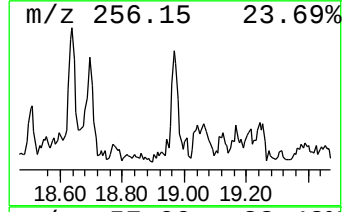
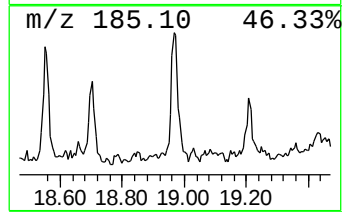
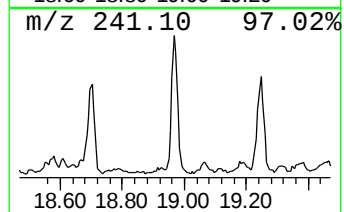
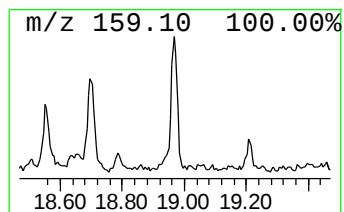
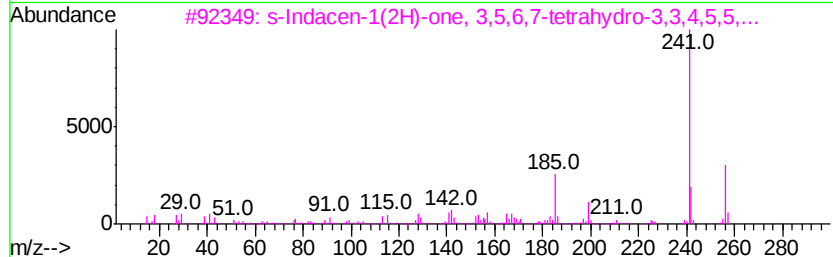
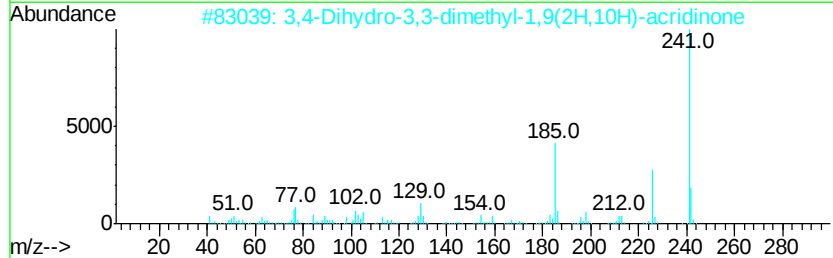
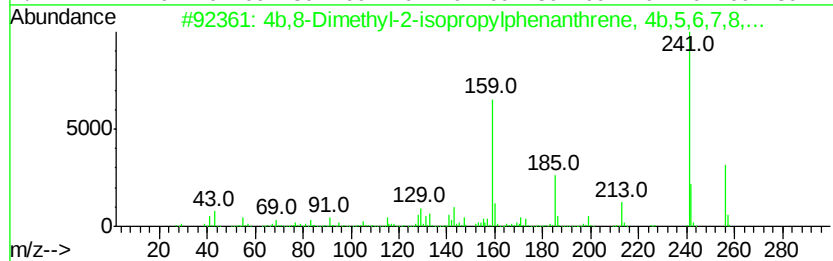
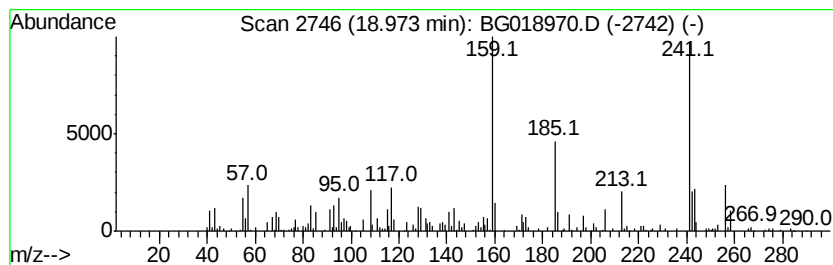
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 48 4b,8-Dimethyl-2-isopropylph... Concentration Rank 69

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.97 | 2.99 ng/ul | 177076 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 4b,8-Dimethyl-2-isopropylphenant... | 256 | C19H28    | 1000197-14-1 | 93   |
| 2    |    | 3,4-Dihydro-3,3-dimethyl-1,9(2H,... | 241 | C15H15N02 | 1000212-95-2 | 38   |
| 3    |    | s-Indacen-1(2H)-one, 3,5,6,7-tet... | 256 | C18H24O   | 038754-94-8  | 38   |
| 4    |    | 4-Ethyl-9-chloro-acridine           | 241 | C15H12ClN | 065753-76-6  | 25   |
| 5    |    | Benzene, 1,1'-(1-methylethyliden... | 256 | C17H20O2  | 001568-83-8  | 22   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

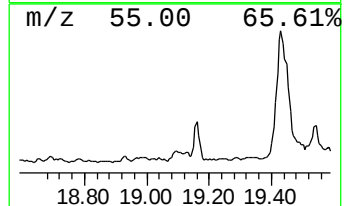
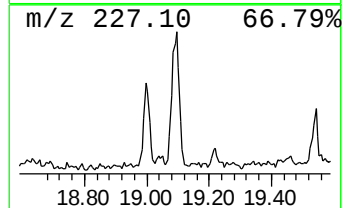
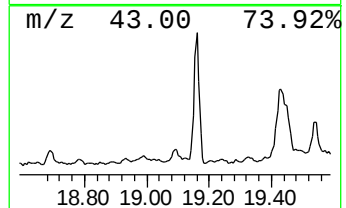
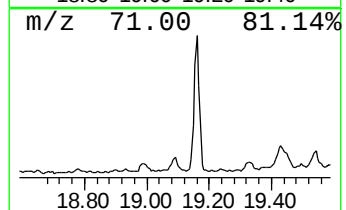
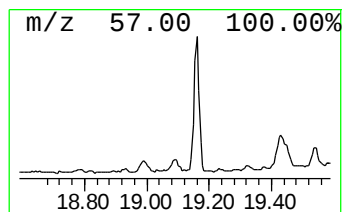
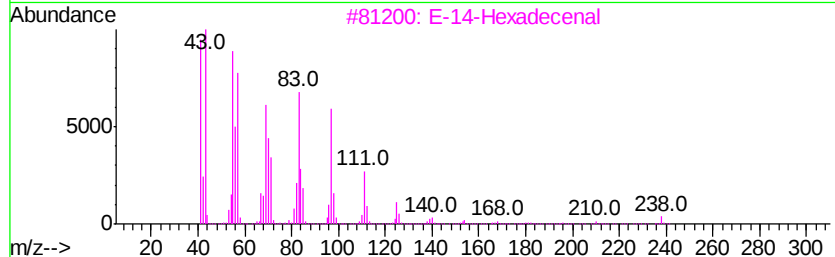
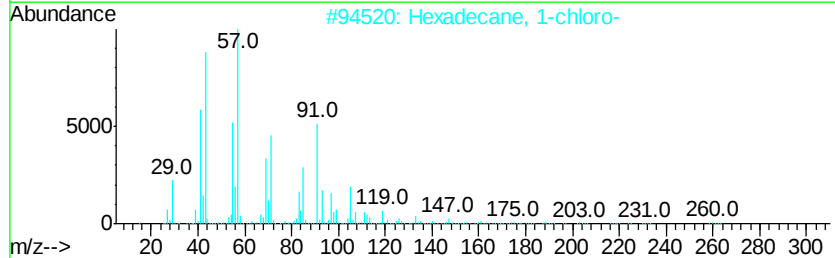
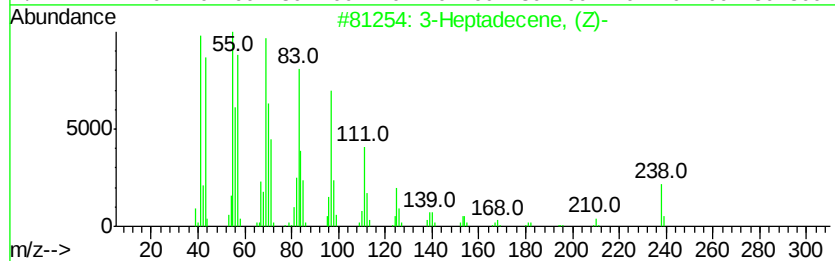
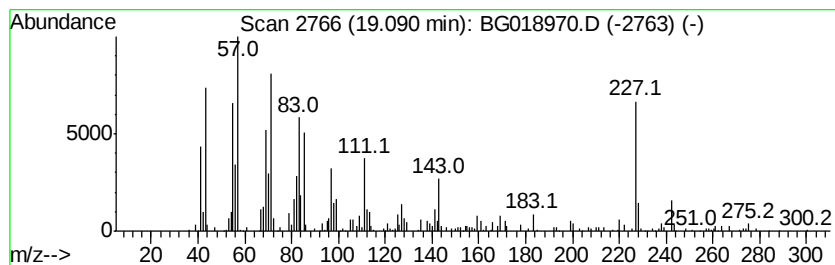
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 49 (DEL) Alkane: Cyclic19.09 Concentration Rank 64

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.09 | 3.33 ng/ul | 197221 | Phenanthrene-d10 | 17.37 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm  | CAS#         | Qual |
|------|----|---|---------------------------|-----|----------|--------------|------|
| 1    |    |   | 3-Heptadecene, (Z)-       | 238 | C17H34   | 1000141-67-3 | 66   |
| 2    |    |   | Hexadecane, 1-chloro-     | 260 | C16H33Cl | 004860-03-1  | 53   |
| 3    |    |   | E-14-Hexadecenal          | 238 | C16H30O  | 330207-53-9  | 45   |
| 4    |    |   | 1-Hexacosanol             | 382 | C26H54O  | 000506-52-5  | 38   |
| 5    |    |   | Octane, 5-ethyl-2-methyl- | 156 | C11H24   | 062016-18-6  | 38   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

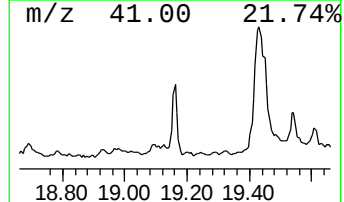
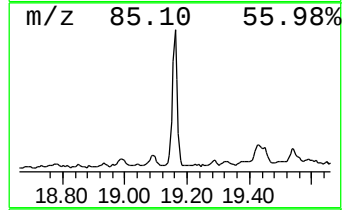
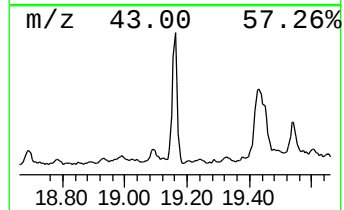
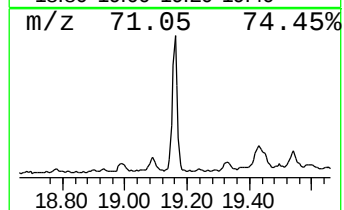
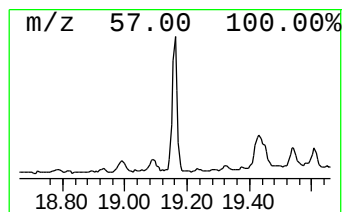
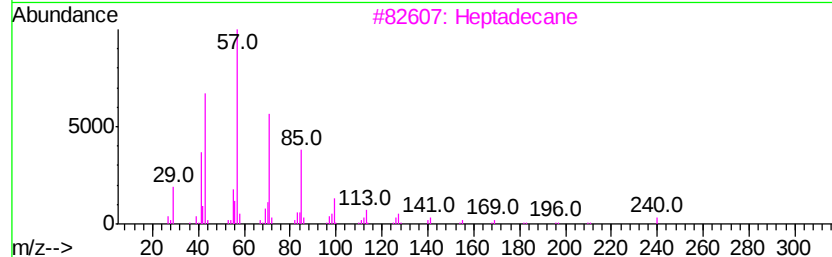
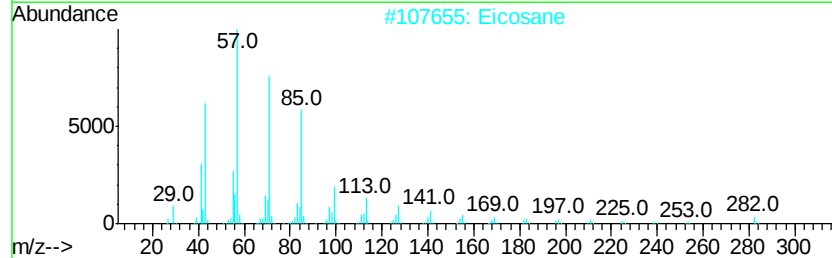
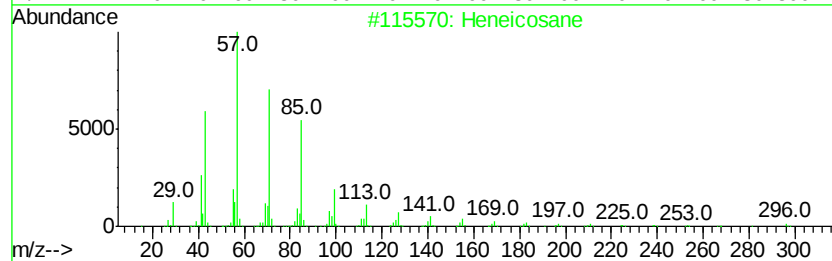
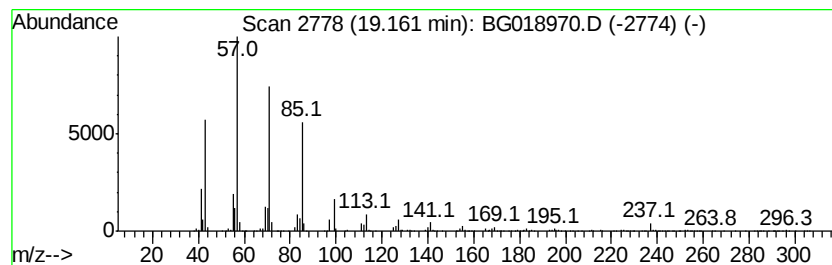
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 19.16 | 16.79 ng/ul | 995141 | Phenanthrene-d10 | 17.37 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane  | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    | Eicosane     | 282 | C20H42  | 000112-95-8 | 91   |
| 3    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 90   |
| 4    |    | Eicosane     | 282 | C20H42  | 000112-95-8 | 90   |
| 5    |    | Nonadecane   | 268 | C19H40  | 000629-92-5 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

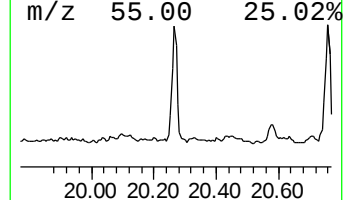
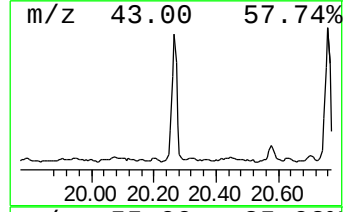
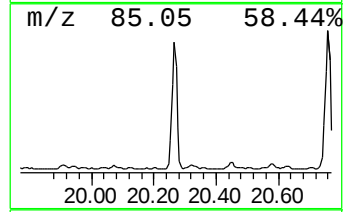
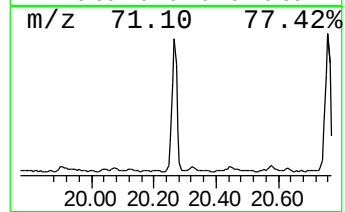
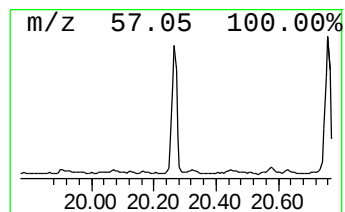
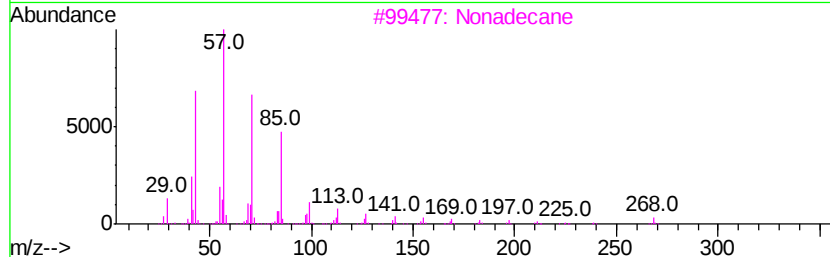
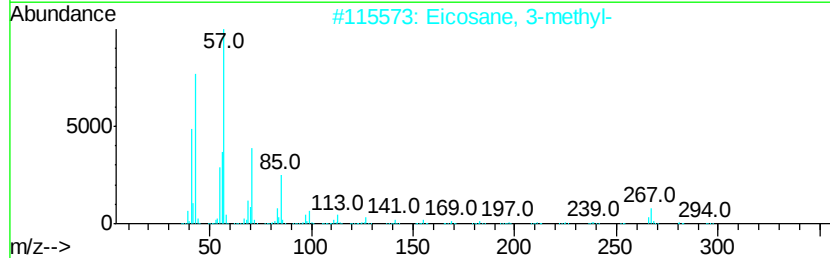
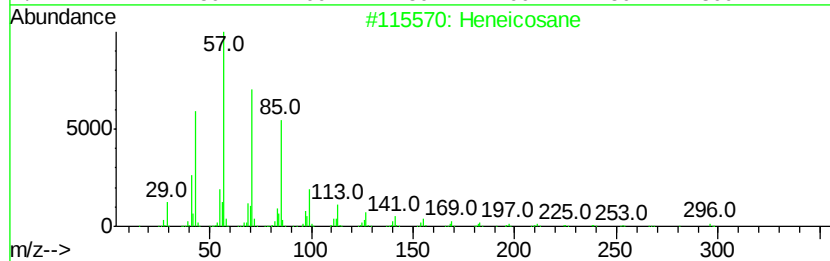
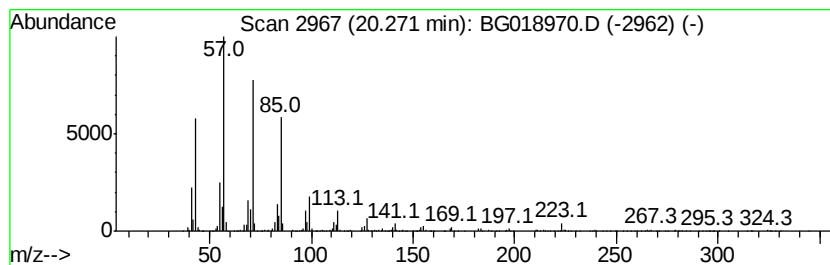
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 20.27 | 23.39 ng/ul | 2023360 | Chrysene-d12     | 21.62 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane            | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    | Eicosane, 3-methyl-    | 296 | C21H44  | 006418-46-8 | 91   |
| 3    |    | Nonadecane             | 268 | C19H40  | 000629-92-5 | 90   |
| 4    |    | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 90   |
| 5    |    | Eicosane               | 282 | C20H42  | 000112-95-8 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

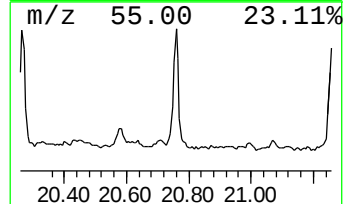
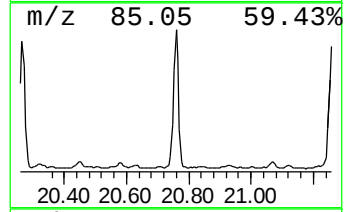
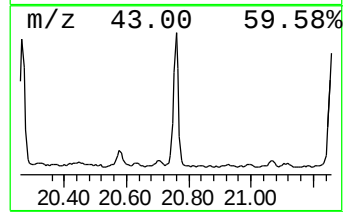
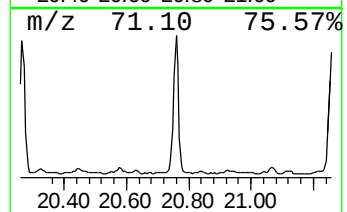
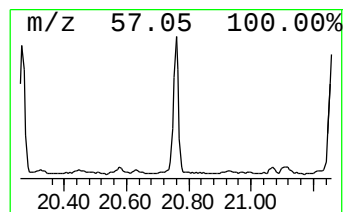
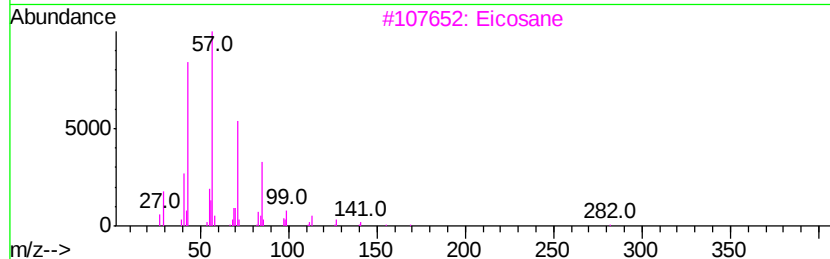
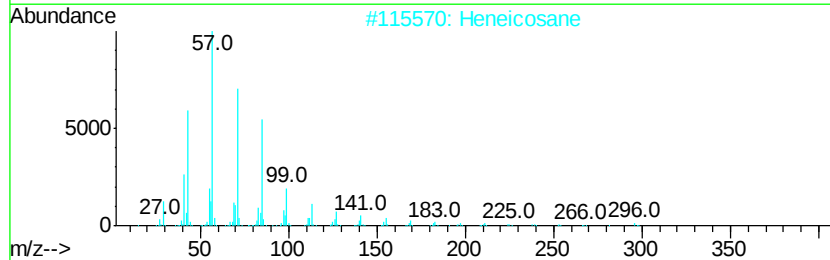
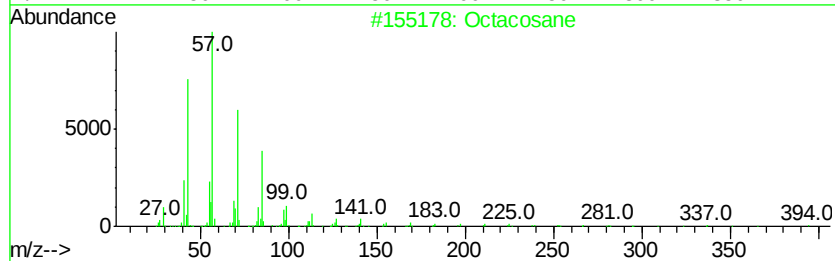
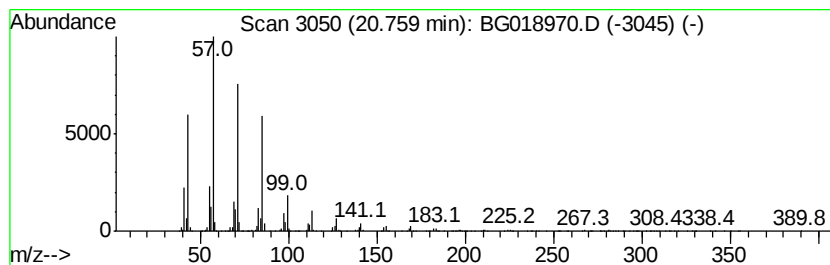
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 20.76 | 26.05 ng/ul | 2253110 | Chrysene-d12     | 21.62 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Octacosane   | 394 | C28H58  | 000630-02-4 | 91   |
| 2    |    | Heneicosane  | 296 | C21H44  | 000629-94-7 | 90   |
| 3    |    | Eicosane     | 282 | C20H42  | 000112-95-8 | 90   |
| 4    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 87   |
| 5    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

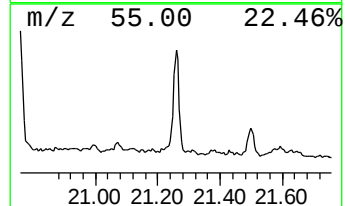
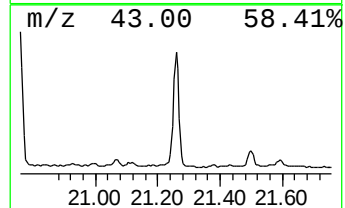
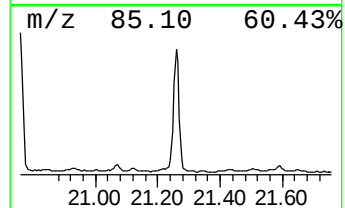
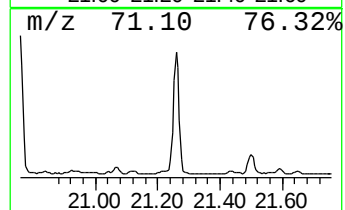
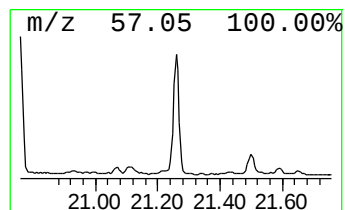
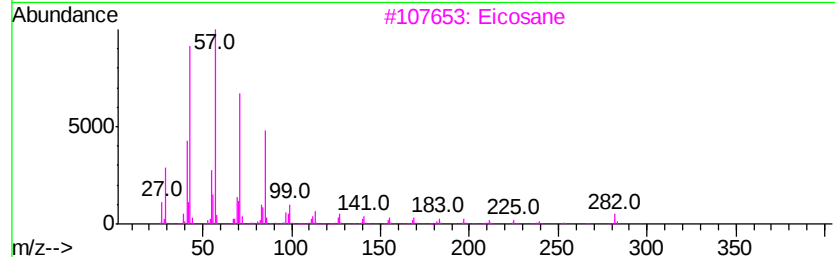
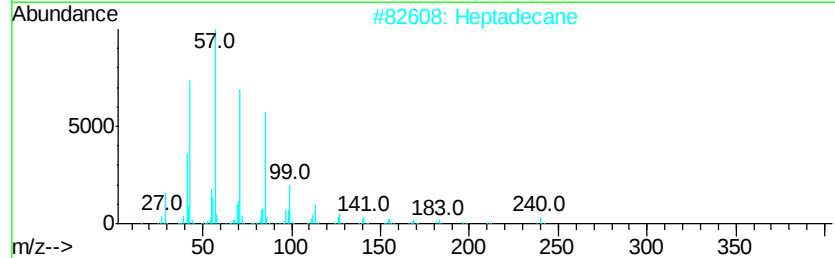
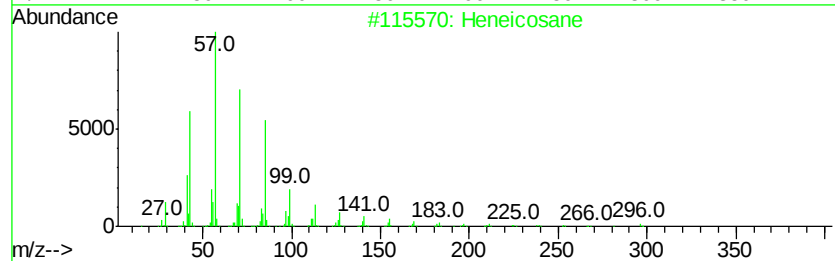
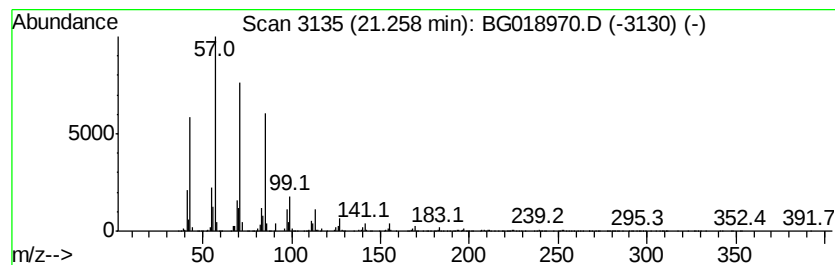
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 21.26 | 25.97 ng/ul | 2246500 | Chrysene-d12     | 21.62 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane            | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    | Heptadecane            | 240 | C17H36  | 000629-78-7 | 91   |
| 3    |    | Eicosane               | 282 | C20H42  | 000112-95-8 | 91   |
| 4    |    | Octacosane             | 394 | C28H58  | 000630-02-4 | 91   |
| 5    |    | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

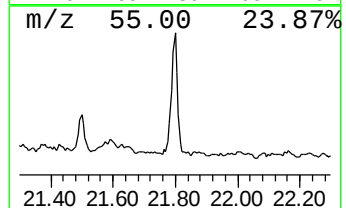
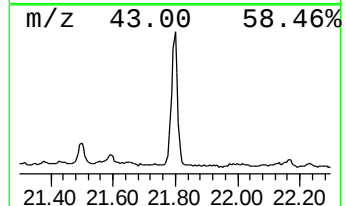
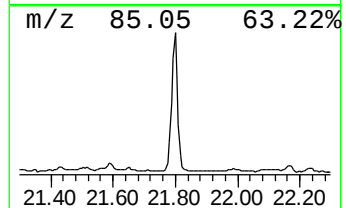
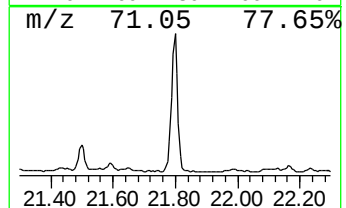
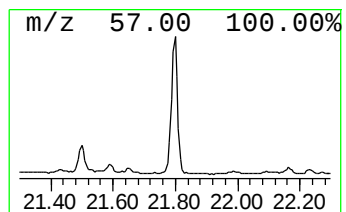
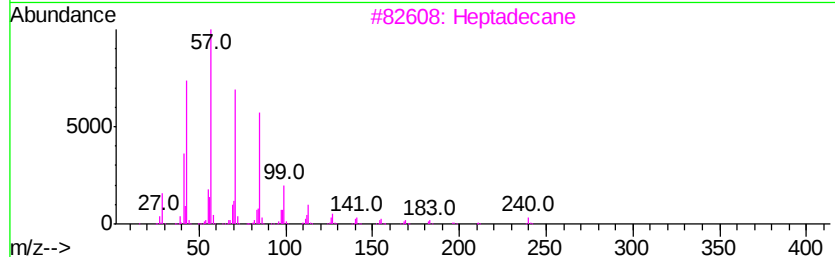
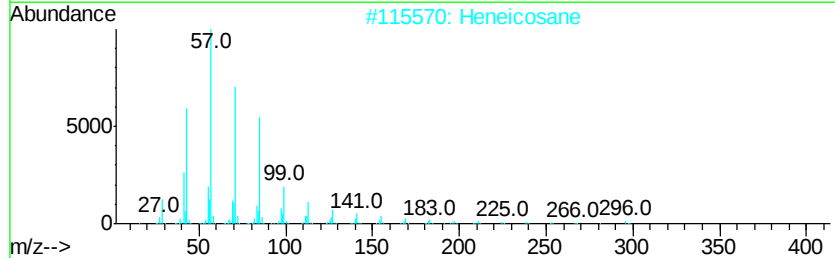
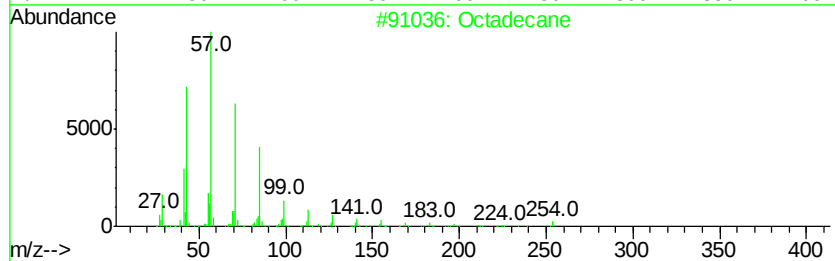
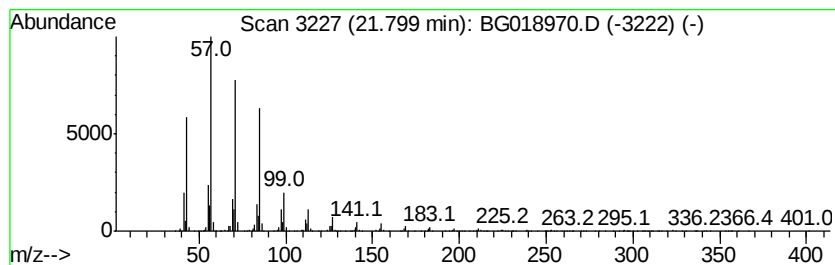
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 64 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 21.80 | 20.84 ng/ul | 1802570 | Chrysene-d12     | 21.62 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Octadecane            | 254 | C18H38  | 000593-45-3 | 96   |
| 2    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    | Heptadecane           | 240 | C17H36  | 000629-78-7 | 91   |
| 4    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |
| 5    |    | Octacosane            | 394 | C28H58  | 000630-02-4 | 90   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

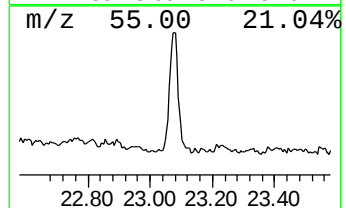
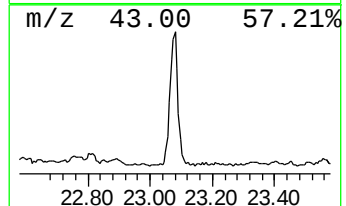
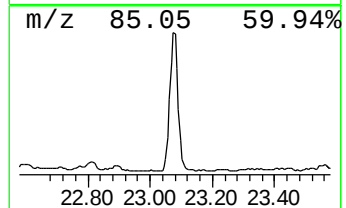
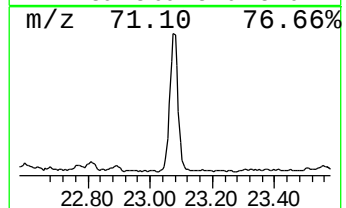
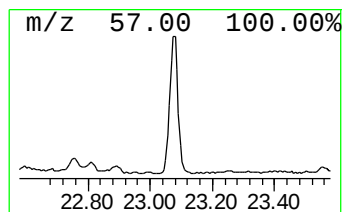
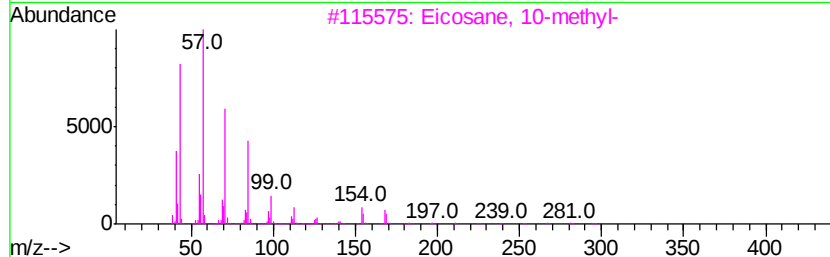
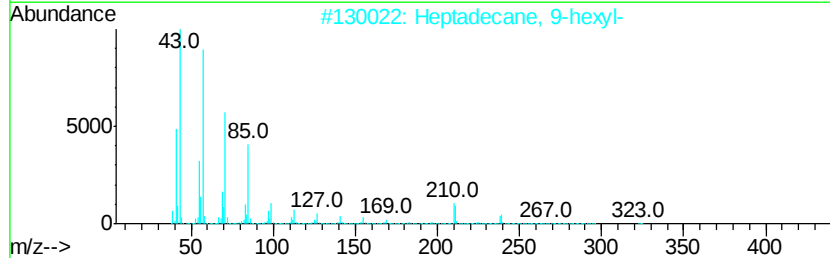
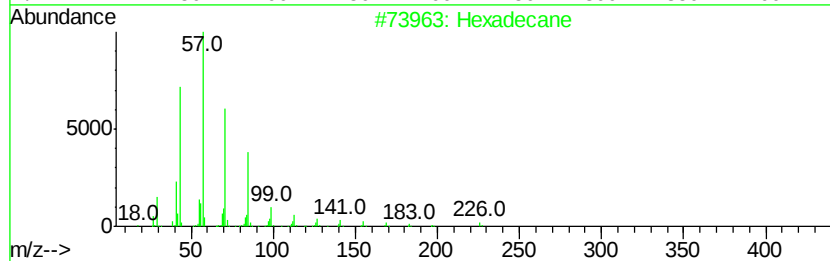
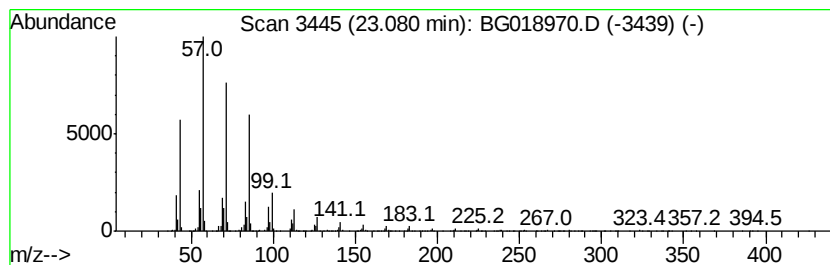
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 68 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 23.08 | 15.71 ng/ul | 1359260 | Chrysene-d12     | 21.62 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane              | 226 | C16H34  | 000544-76-3 | 94   |
| 2    |    |   | Heptadecane, 9-hexyl-   | 324 | C23H48  | 055124-79-3 | 93   |
| 3    |    |   | Eicosane, 10-methyl-    | 296 | C21H44  | 054833-23-7 | 93   |
| 4    |    |   | Heneicosane, 10-methyl- | 310 | C22H46  | 013287-25-7 | 91   |
| 5    |    |   | Heneicosane             | 296 | C21H44  | 000629-94-7 | 91   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

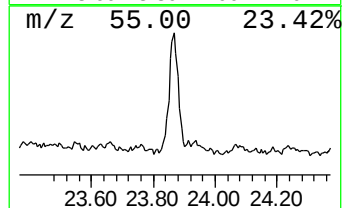
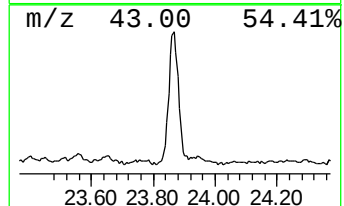
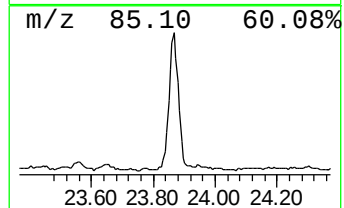
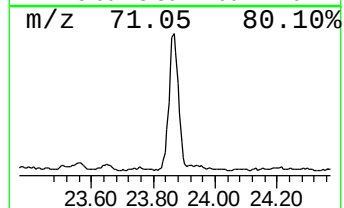
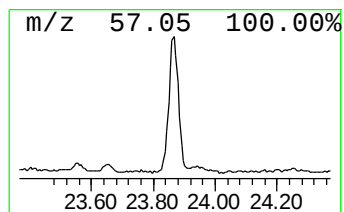
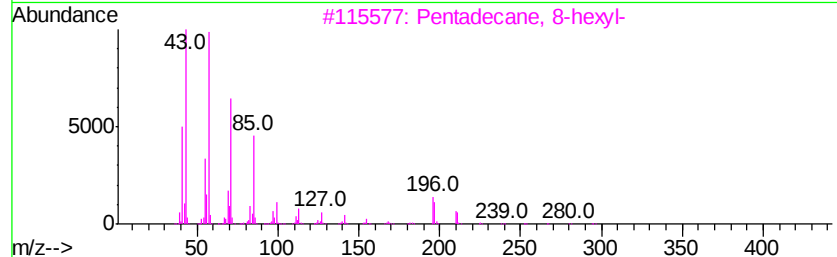
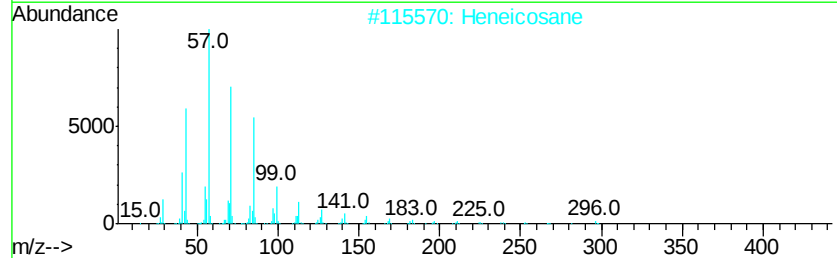
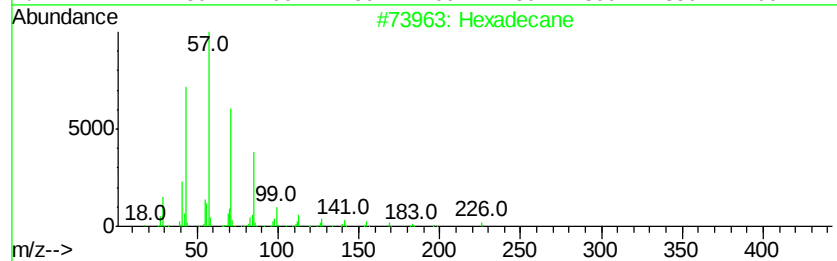
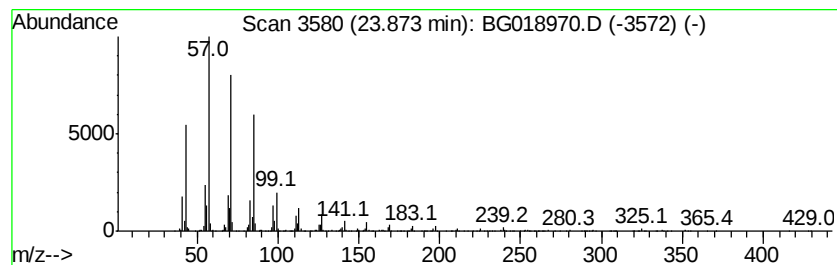
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 69 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 23.87 | 10.09 ng/ul | 1157100 | Perylene-d12     | 24.81 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane            | 226 | C16H34  | 000544-76-3 | 93   |
| 2    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 90   |
| 4    |    | Hexadecane            | 226 | C16H34  | 000544-76-3 | 90   |
| 5    |    | Heptadecane           | 240 | C17H36  | 000629-78-7 | 86   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\  
Data File : BG018970.D  
Acq On : 30 Sep 2015 13:28  
Operator : UM/NP  
Sample : G3690-01DL 5X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M  
Quant Title : SVOA CALIBRATION

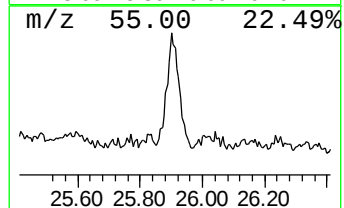
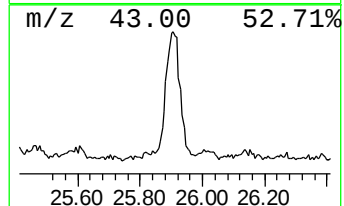
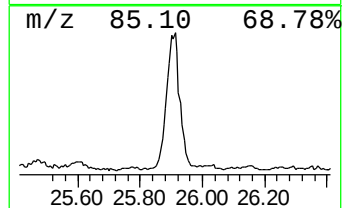
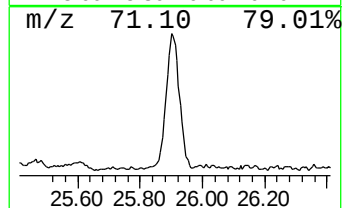
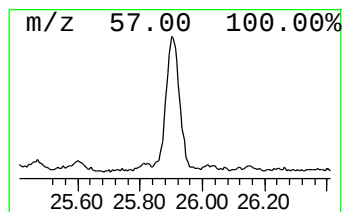
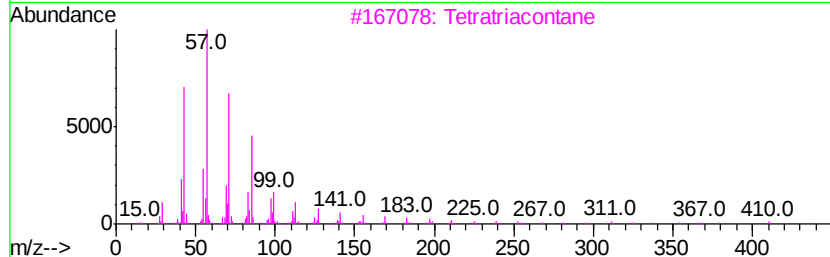
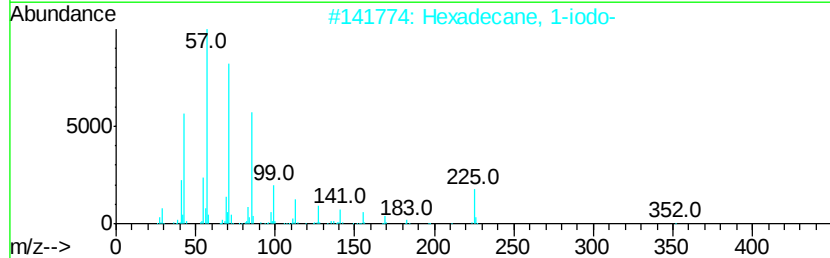
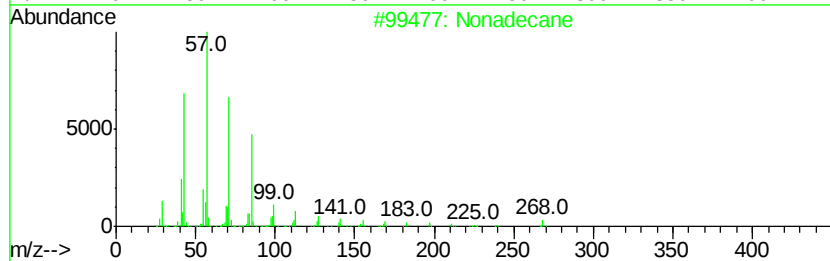
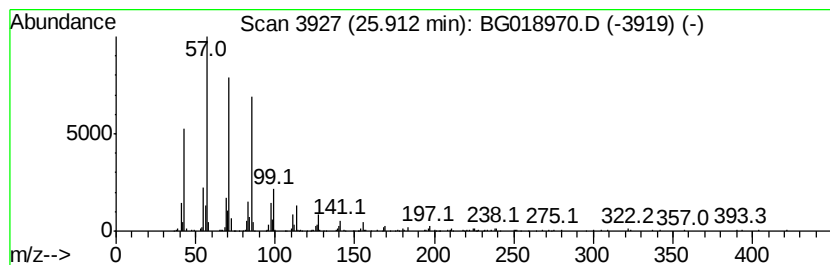
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 49

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 25.91 | 5.33 ng/ul | 611071 | Perylene-d12     | 24.81 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane          | 268 | C19H40  | 000629-92-5 | 93   |
| 2    |    |   | Hexadecane, 1-iodo- | 352 | C16H33I | 000544-77-4 | 93   |
| 3    |    |   | Tetratriacontane    | 479 | C34H70  | 014167-59-0 | 91   |
| 4    |    |   | Octacosane          | 394 | C28H58  | 000630-02-4 | 91   |
| 5    |    |   | Heneicosane         | 296 | C21H44  | 000629-94-7 | 90   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG100115\  
 Data File : BG018970.D  
 Acq On : 30 Sep 2015 13:28  
 Operator : UM/NP  
 Sample : G3690-01DL 5X  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BC2W1DL

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG091015.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 |
| Butane, 2-methoxy... | 3.16  | 15.6    | ng/ul | 590974   | 1                     | 7.99  | 758717  | 20.0 |
| Benzene, 1,3-dime... | 5.62  | 12.1    | ng/ul | 459035   | 1                     | 7.99  | 758717  | 20.0 |
| (DEL) Alkane: Str... | 6.97  | 6.4     | ng/ul | 243552   | 1                     | 7.99  | 758717  | 20.0 |
| (DEL) Alkane: Str... | 7.60  | 16.6    | ng/ul | 627929   | 1                     | 7.99  | 758717  | 20.0 |
| (DEL) Alkane: Cyc... | 7.91  | 4.2     | ng/ul | 160167   | 1                     | 7.99  | 758717  | 20.0 |
| unknown-01           | 8.11  | 7.7     | ng/ul | 291790   | 1                     | 7.99  | 758717  | 20.0 |
| (DEL) Alkane: Str... | 8.21  | 19.3    | ng/ul | 730274   | 1                     | 7.99  | 758717  | 20.0 |
| (DEL) Alkane: Cyc... | 8.27  | 4.6     | ng/ul | 173809   | 1                     | 7.99  | 758717  | 20.0 |
| (DEL) Alkane: Str... | 8.53  | 17.1    | ng/ul | 646849   | 1                     | 7.99  | 758717  | 20.0 |
| (DEL) Alkane: Str... | 8.86  | 4.4     | ng/ul | 166921   | 1                     | 7.99  | 758717  | 20.0 |
| Benzene, 1-methyl... | 9.09  | 6.8     | ng/ul | 257686   | 1                     | 7.99  | 758717  | 20.0 |
| (DEL) Alkane: Str... | 9.21  | 26.4    | ng/ul | 1001710  | 1                     | 7.99  | 758717  | 20.0 |
| unknown-02           | 9.34  | 5.1     | ng/ul | 192788   | 1                     | 7.99  | 758717  | 20.0 |
| Naphthalene, deca... | 9.97  | 5.4     | ng/ul | 147348   | 2                     | 10.79 | 541673  | 20.0 |
| (DEL) Alkane: Str... | 10.19 | 7.1     | ng/ul | 192316   | 2                     | 10.79 | 541673  | 20.0 |
| Parachlorophenol     | 10.66 | 35.2    | ng/ul | 952422   | 2                     | 10.79 | 541673  | 20.0 |
| Phenol, p-tert-bu... | 12.12 | 7.4     | ng/ul | 199577   | 2                     | 10.79 | 541673  | 20.0 |
| (DEL) Alkane: Str... | 12.19 | 6.3     | ng/ul | 169942   | 2                     | 10.79 | 541673  | 20.0 |
| Phenol, 3,4-dichl... | 13.64 | 3.4     | ng/ul | 174155   | 3                     | 14.61 | 1029310 | 20.0 |
| Diphenyl ether       | 13.66 | 3.5     | ng/ul | 178469   | 3                     | 14.61 | 1029310 | 20.0 |
| 1-Undecanol          | 14.25 | 3.1     | ng/ul | 161716   | 3                     | 14.61 | 1029310 | 20.0 |
| (DEL) Alkane: Cyc... | 14.37 | 4.0     | ng/ul | 207050   | 3                     | 14.61 | 1029310 | 20.0 |
| (DEL) Alkane: Str... | 14.50 | 3.4     | ng/ul | 176861   | 3                     | 14.61 | 1029310 | 20.0 |
| 1,1'-Biphenyl, 4-... | 14.70 | 2.7     | ng/ul | 139796   | 3                     | 14.61 | 1029310 | 20.0 |
| o-Hydroxybiphenyl    | 14.90 | 124.1   | ng/ul | 6388840  | 3                     | 14.61 | 1029310 | 20.0 |
| Dodecanoic acid      | 15.04 | 3.5     | ng/ul | 177354   | 3                     | 14.61 | 1029310 | 20.0 |
| (DEL) Alkane: Str... | 15.45 | 2.9     | ng/ul | 147098   | 3                     | 14.61 | 1029310 | 20.0 |
| (DEL) Alkane: Cyc... | 16.15 | 5.5     | ng/ul | 326583   | 4                     | 17.37 | 1185270 | 20.0 |
| (DEL) Alkane: Str... | 16.31 | 7.0     | ng/ul | 413367   | 4                     | 17.37 | 1185270 | 20.0 |
| 1-[1-[2-Pyridyl]e... | 16.34 | 6.2     | ng/ul | 368702   | 4                     | 17.37 | 1185270 | 20.0 |
| Phenol, 4-(1,1,3,... | 16.43 | 3.3     | ng/ul | 196154   | 4                     | 17.37 | 1185270 | 20.0 |
| Acetamide, N-(3-m... | 16.49 | 3.6     | ng/ul | 216255   | 4                     | 17.37 | 1185270 | 20.0 |
| unknown-03           | 16.55 | 4.9     | ng/ul | 288350   | 4                     | 17.37 | 1185270 | 20.0 |
| unknown-04           | 16.76 | 5.7     | ng/ul | 340418   | 4                     | 17.37 | 1185270 | 20.0 |
| (DEL) Alkane: Str... | 17.16 | 4.5     | ng/ul | 265043   | 4                     | 17.37 | 1185270 | 20.0 |
| 1-(4-Fluorophenyl... | 17.60 | 2.3     | ng/ul | 134570   | 4                     | 17.37 | 1185270 | 20.0 |
| 1,3-Dioxolane, 2-... | 17.71 | 26.2    | ng/ul | 1553990  | 4                     | 17.37 | 1185270 | 20.0 |
| Butyric acid, 3-m... | 17.76 | 3.1     | ng/ul | 181591   | 4                     | 17.37 | 1185270 | 20.0 |
| (DEL) Alkane: Str... | 17.84 | 12.1    | ng/ul | 718348   | 4                     | 17.37 | 1185270 | 20.0 |
| Clorophene           | 18.06 | 6.1     | ng/ul | 359732   | 4                     | 17.37 | 1185270 | 20.0 |
| Benzene, 1,1',1'...  | 18.44 | 10.5    | ng/ul | 621461   | 4                     | 17.37 | 1185270 | 20.0 |
| Benzoic acid, 4-(... | 18.55 | 17.9    | ng/ul | 1058170  | 4                     | 17.37 | 1185270 | 20.0 |
| unknown-05           | 18.66 | 2.3     | ng/ul | 136748   | 4                     | 17.37 | 1185270 | 20.0 |
| 9,10-Anthracenedi... | 18.69 | 12.7    | ng/ul | 755061   | 4                     | 17.37 | 1185270 | 20.0 |
| unknown-06           | 18.78 | 3.1     | ng/ul | 181414   | 4                     | 17.37 | 1185270 | 20.0 |
| Phenanthrene, 7-e... | 18.93 | 3.6     | ng/ul | 212388   | 4                     | 17.37 | 1185270 | 20.0 |
| 4b,8-Dimethyl-2-i... | 18.97 | 3.0     | ng/ul | 177076   | 4                     | 17.37 | 1185270 | 20.0 |
| (DEL) Alkane: Cyc... | 19.09 | 3.3     | ng/ul | 197221   | 4                     | 17.37 | 1185270 | 20.0 |
| (DEL) Alkane: Str... | 19.16 | 16.8    | ng/ul | 995141   | 4                     | 17.37 | 1185270 | 20.0 |

|       |                |       |      |       |         |   |       |         |      |
|-------|----------------|-------|------|-------|---------|---|-------|---------|------|
| (DEL) | Alkane: Str... | 20.27 | 23.4 | ng/ul | 2023360 | 5 | 21.62 | 1730010 | 20.0 |
| (DEL) | Alkane: Str... | 20.76 | 26.1 | ng/ul | 2253110 | 5 | 21.62 | 1730010 | 20.0 |
| (DEL) | Alkane: Str... | 21.26 | 26.0 | ng/ul | 2246500 | 5 | 21.62 | 1730010 | 20.0 |
| (DEL) | Alkane: Str... | 21.80 | 20.8 | ng/ul | 1802570 | 5 | 21.62 | 1730010 | 20.0 |
| (DEL) | Alkane: Str... | 23.08 | 15.7 | ng/ul | 1359260 | 5 | 21.62 | 1730010 | 20.0 |
| (DEL) | Alkane: Str... | 23.87 | 10.1 | ng/ul | 1157100 | 6 | 24.81 | 2294260 | 20.0 |
| (DEL) | Alkane: Str... | 25.91 | 5.3  | ng/ul | 611071  | 6 | 24.81 | 2294260 | 20.0 |

Instrument :  
BNA\_G  
ClientSampleId :  
BC2W1DL