

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
 Data File : BG018305.D  
 Acq On : 6 Aug 2015 15:20  
 Operator : UM/TP  
 Sample : G3109-15DL 5X  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C01W3DL

## 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-BG080515.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 2.811       | 17            | 21          | 23           | rVB4     | 3337           | 3029          | 0.31%           | 0.030%        |
| 2         | 2.923       | 38            | 40          | 42           | rVB2     | 3115           | 2497          | 0.26%           | 0.025%        |
| 3         | 3.040       | 57            | 60          | 62           | rBV2     | 1875           | 2307          | 0.24%           | 0.023%        |
| 4         | 3.117       | 72            | 73          | 76           | rBV2     | 3483           | 2542          | 0.26%           | 0.025%        |
| 5         | 3.240       | 90            | 94          | 98           | rVB3     | 5380           | 7293          | 0.75%           | 0.072%        |
| 6         | 3.428       | 122           | 126         | 129          | rBV4     | 3281           | 5089          | 0.52%           | 0.050%        |
| 7         | 4.527       | 309           | 313         | 316          | rVB2     | 2125           | 2932          | 0.30%           | 0.029%        |
| 8         | 5.438       | 466           | 468         | 471          | rBV2     | 2807           | 3352          | 0.34%           | 0.033%        |
| 9         | 5.631       | 498           | 501         | 503          | rBV2     | 2281           | 2466          | 0.25%           | 0.024%        |
| 10        | 5.778       | 523           | 526         | 529          | rBV2     | 1666           | 2436          | 0.25%           | 0.024%        |
| 11        | 6.642       | 668           | 673         | 681          | rBV      | 16441          | 25473         | 2.61%           | 0.251%        |
| 12        | 6.771       | 690           | 695         | 704          | rVB2     | 11123          | 19049         | 1.95%           | 0.188%        |
| 13        | 6.971       | 723           | 729         | 736          | rVB3     | 15303          | 26603         | 2.73%           | 0.263%        |
| 14        | 7.071       | 744           | 746         | 751          | rVB      | 2672           | 3785          | 0.39%           | 0.037%        |
| 15        | 7.341       | 790           | 792         | 795          | rBV      | 2382           | 2785          | 0.29%           | 0.027%        |
| 16        | 7.429       | 801           | 807         | 815          | rVB      | 99294          | 149949        | 15.36%          | 1.480%        |
| 17        | 7.788       | 863           | 868         | 870          | rBV      | 1681           | 2684          | 0.27%           | 0.026%        |
| 18        | 7.987       | 901           | 902         | 905          | rVB      | 3262           | 2536          | 0.26%           | 0.025%        |
| 19        | 8.023       | 905           | 908         | 909          | rBV      | 2585           | 2385          | 0.24%           | 0.024%        |
| 20        | 8.070       | 912           | 916         | 917          | rBV      | 3014           | 3416          | 0.35%           | 0.034%        |
| 21        | 8.170       | 927           | 933         | 939          | rVB4     | 12568          | 21716         | 2.22%           | 0.214%        |
| 22        | 8.587       | 997           | 1004        | 1009         | rVB2     | 15453          | 25202         | 2.58%           | 0.249%        |
| 23        | 8.851       | 1044          | 1049        | 1050         | rBV      | 1914           | 2503          | 0.26%           | 0.025%        |
| 24        | 8.939       | 1060          | 1064        | 1066         | rVB      | 2068           | 2939          | 0.30%           | 0.029%        |
| 25        | 9.174       | 1101          | 1104        | 1107         | rBV      | 2006           | 2412          | 0.25%           | 0.024%        |
| 26        | 9.304       | 1120          | 1126        | 1131         | rBV      | 16315          | 23812         | 2.44%           | 0.235%        |
| 27        | 9.862       | 1215          | 1221        | 1229         | rBV2     | 21275          | 34711         | 3.56%           | 0.343%        |
| 28        | 10.214      | 1273          | 1281        | 1286         | rBV      | 130888         | 218055        | 22.34%          | 2.153%        |
| 29        | 10.267      | 1286          | 1290        | 1296         | rVB2     | 14530          | 23157         | 2.37%           | 0.229%        |
| 30        | 11.054      | 1421          | 1424        | 1427         | rVV2     | 2183           | 2343          | 0.24%           | 0.023%        |
| 31        | 11.895      | 1559          | 1567        | 1574         | rBV      | 16444          | 25009         | 2.56%           | 0.247%        |
| 32        | 12.012      | 1585          | 1587        | 1589         | rBV      | 2760           | 2850          | 0.29%           | 0.028%        |
| 33        | 12.118      | 1601          | 1605        | 1606         | rBV3     | 8693           | 10600         | 1.09%           | 0.105%        |
| 34        | 12.647      | 1692          | 1695        | 1699         | rVB2     | 2933           | 4184          | 0.43%           | 0.041%        |

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 Max Peaks: 100  
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Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG080515.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 12.935 | 1740 | 1744 | 1749 | rVB3 | 6638   | 8934   | 0.92%  | 0.088% |
| 36 | 13.129 | 1773 | 1777 | 1784 | rVV4 | 4300   | 10557  | 1.08%  | 0.104% |
| 37 | 13.270 | 1797 | 1801 | 1802 | rBV3 | 7299   | 8184   | 0.84%  | 0.081% |
| 38 | 13.287 | 1802 | 1804 | 1809 | rVB2 | 9248   | 11033  | 1.13%  | 0.109% |
| 39 | 13.428 | 1823 | 1828 | 1834 | rVV2 | 12661  | 24478  | 2.51%  | 0.242% |
| 40 | 13.487 | 1834 | 1838 | 1841 | rVV3 | 8706   | 14297  | 1.46%  | 0.141% |
| 41 | 13.528 | 1841 | 1845 | 1850 | rVV2 | 45763  | 64109  | 6.57%  | 0.633% |
| 42 | 13.575 | 1850 | 1853 | 1856 | rVV2 | 18163  | 24634  | 2.52%  | 0.243% |
| 43 | 13.604 | 1856 | 1858 | 1864 | rVV4 | 6956   | 8624   | 0.88%  | 0.085% |
| 44 | 13.663 | 1864 | 1868 | 1873 | rVV4 | 7973   | 12588  | 1.29%  | 0.124% |
| 45 | 13.698 | 1873 | 1874 | 1878 | rVB2 | 3188   | 3147   | 0.32%  | 0.031% |
| 46 | 13.804 | 1885 | 1892 | 1896 | rBV2 | 42392  | 66402  | 6.80%  | 0.656% |
| 47 | 14.033 | 1926 | 1931 | 1936 | rBV4 | 3775   | 6786   | 0.70%  | 0.067% |
| 48 | 14.116 | 1936 | 1945 | 1951 | rBV2 | 232806 | 348759 | 35.73% | 3.443% |
| 49 | 14.174 | 1951 | 1955 | 1966 | rVB  | 61167  | 92737  | 9.50%  | 0.916% |
| 50 | 14.257 | 1966 | 1969 | 1972 | rVB3 | 3636   | 3444   | 0.35%  | 0.034% |
| 51 | 14.380 | 1988 | 1990 | 1993 | rBV4 | 5683   | 7620   | 0.78%  | 0.075% |
| 52 | 14.515 | 2008 | 2013 | 2019 | rVB  | 42414  | 66404  | 6.80%  | 0.656% |
| 53 | 14.597 | 2023 | 2027 | 2032 | rBV7 | 6256   | 10610  | 1.09%  | 0.105% |
| 54 | 14.656 | 2035 | 2037 | 2039 | rBV3 | 4707   | 4469   | 0.46%  | 0.044% |
| 55 | 14.744 | 2048 | 2052 | 2056 | rBV3 | 6702   | 11361  | 1.16%  | 0.112% |
| 56 | 14.791 | 2058 | 2060 | 2064 | rVB3 | 7967   | 9772   | 1.00%  | 0.096% |
| 57 | 14.897 | 2075 | 2078 | 2081 | rBV4 | 6708   | 10869  | 1.11%  | 0.107% |
| 58 | 15.114 | 2111 | 2115 | 2120 | rVB  | 52976  | 74763  | 7.66%  | 0.738% |
| 59 | 15.173 | 2120 | 2125 | 2129 | rVB  | 96213  | 135076 | 13.84% | 1.334% |
| 60 | 15.208 | 2129 | 2131 | 2134 | rBV3 | 5369   | 7629   | 0.78%  | 0.075% |
| 61 | 15.267 | 2137 | 2141 | 2143 | rBV3 | 19034  | 24389  | 2.50%  | 0.241% |
| 62 | 15.297 | 2144 | 2146 | 2149 | rVB2 | 17184  | 15552  | 1.59%  | 0.154% |
| 63 | 15.332 | 2149 | 2152 | 2156 | rVB2 | 17298  | 21627  | 2.22%  | 0.214% |
| 64 | 15.385 | 2158 | 2161 | 2165 | rBV6 | 12342  | 22026  | 2.26%  | 0.217% |
| 65 | 15.467 | 2173 | 2175 | 2184 | rVB  | 23878  | 39705  | 4.07%  | 0.392% |
| 66 | 15.620 | 2197 | 2201 | 2209 | rVB2 | 24578  | 44509  | 4.56%  | 0.439% |
| 67 | 15.972 | 2256 | 2261 | 2264 | rBV4 | 16448  | 26170  | 2.68%  | 0.258% |
| 68 | 16.037 | 2268 | 2272 | 2278 | rVB8 | 14500  | 24873  | 2.55%  | 0.246% |
| 69 | 16.231 | 2302 | 2305 | 2311 | rVB5 | 20125  | 29678  | 3.04%  | 0.293% |
| 70 | 16.301 | 2311 | 2317 | 2319 | rBV7 | 12759  | 23988  | 2.46%  | 0.237% |
| 71 | 16.525 | 2353 | 2355 | 2360 | rVB6 | 10987  | 12941  | 1.33%  | 0.128% |

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Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG080515.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |        |        |         |        |
|-----|--------|------|------|------|------|--------|--------|---------|--------|
| 72  | 16.695 | 2380 | 2384 | 2388 | rVB  | 57199  | 68756  | 7.04%   | 0.679% |
| 73  | 16.877 | 2409 | 2415 | 2418 | rBV  | 273234 | 408629 | 41.86%  | 4.034% |
| 74  | 16.918 | 2418 | 2422 | 2427 | rVV  | 763106 | 976069 | 100.00% | 9.637% |
| 75  | 16.977 | 2427 | 2432 | 2434 | rVV2 | 66654  | 107719 | 11.04%  | 1.064% |
| 76  | 17.006 | 2434 | 2437 | 2442 | rVB  | 180250 | 227804 | 23.34%  | 2.249% |
| 77  | 17.288 | 2482 | 2485 | 2497 | rVB  | 30907  | 56968  | 5.84%   | 0.562% |
| 78  | 17.394 | 2500 | 2503 | 2506 | rBV4 | 17032  | 18736  | 1.92%   | 0.185% |
| 79  | 17.459 | 2510 | 2514 | 2518 | rBV4 | 24568  | 47846  | 4.90%   | 0.472% |
| 80  | 17.529 | 2522 | 2526 | 2529 | rVB  | 43863  | 54532  | 5.59%   | 0.538% |
| 81  | 17.753 | 2560 | 2564 | 2568 | rVV  | 97867  | 125300 | 12.84%  | 1.237% |
| 82  | 17.800 | 2568 | 2572 | 2578 | rVB  | 141622 | 195768 | 20.06%  | 1.933% |
| 83  | 17.946 | 2592 | 2597 | 2601 | rBV3 | 167128 | 291032 | 29.82%  | 2.873% |
| 84  | 18.170 | 2630 | 2635 | 2640 | rBV4 | 61841  | 102976 | 10.55%  | 1.017% |
| 85  | 18.264 | 2648 | 2651 | 2654 | rVB  | 55454  | 63537  | 6.51%   | 0.627% |
| 86  | 18.505 | 2687 | 2692 | 2698 | rVB  | 564451 | 749306 | 76.77%  | 7.398% |
| 87  | 18.646 | 2712 | 2716 | 2720 | rBV  | 148210 | 173896 | 17.82%  | 1.717% |
| 88  | 18.951 | 2763 | 2768 | 2773 | rBV  | 597911 | 779277 | 79.84%  | 7.694% |
| 89  | 19.010 | 2774 | 2778 | 2783 | rVB  | 491919 | 617145 | 63.23%  | 6.093% |
| 90  | 19.292 | 2821 | 2826 | 2827 | rBV2 | 159313 | 213543 | 21.88%  | 2.108% |
| 91  | 19.321 | 2827 | 2831 | 2840 | rVB2 | 634621 | 893932 | 91.58%  | 8.826% |
| 92  | 19.744 | 2899 | 2903 | 2906 | rVB  | 124601 | 150883 | 15.46%  | 1.490% |
| 93  | 19.921 | 2930 | 2933 | 2937 | rVB  | 77171  | 83189  | 8.52%   | 0.821% |
| 94  | 20.814 | 3083 | 3085 | 3089 | rVB3 | 101968 | 105004 | 10.76%  | 1.037% |
| 95  | 21.084 | 3126 | 3131 | 3135 | rBV2 | 290657 | 520607 | 53.34%  | 5.140% |
| 96  | 21.125 | 3135 | 3138 | 3145 | rVB  | 167403 | 214886 | 22.02%  | 2.122% |
| 97  | 21.248 | 3156 | 3159 | 3163 | rVB2 | 103892 | 123807 | 12.68%  | 1.222% |
| 98  | 21.671 | 3228 | 3231 | 3235 | rVB2 | 146596 | 168091 | 17.22%  | 1.660% |
| 99  | 22.112 | 3304 | 3306 | 3314 | rVB  | 160904 | 205464 | 21.05%  | 2.029% |
| 100 | 22.606 | 3386 | 3390 | 3395 | rBV3 | 242458 | 379142 | 38.84%  | 3.743% |

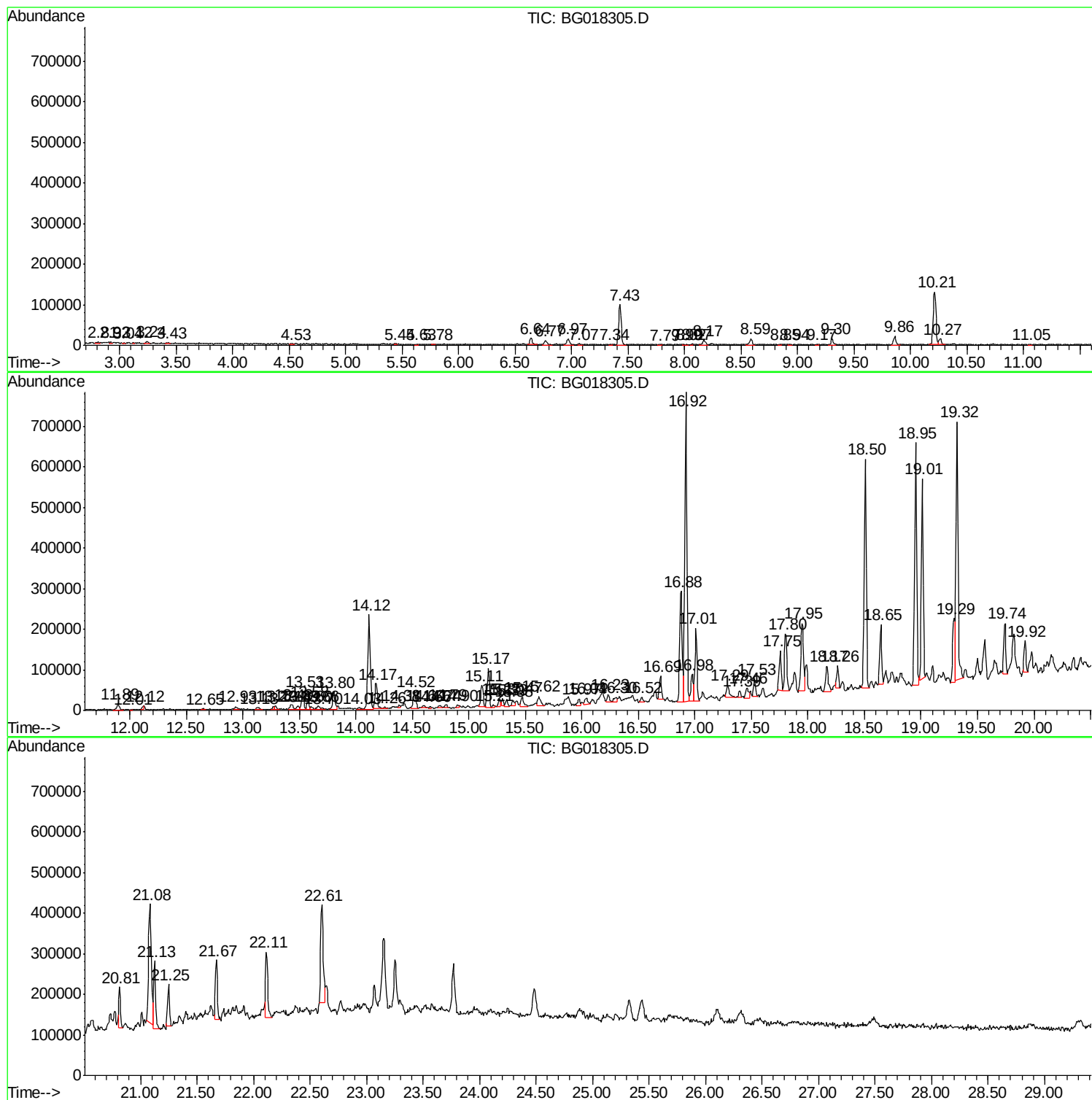
Sum of corrected areas: 10128689

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
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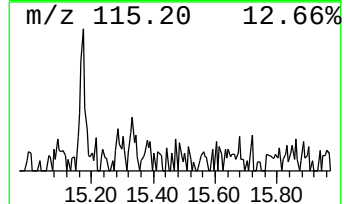
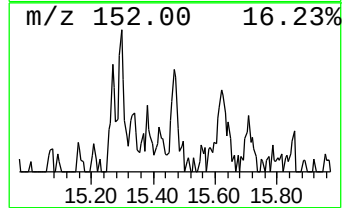
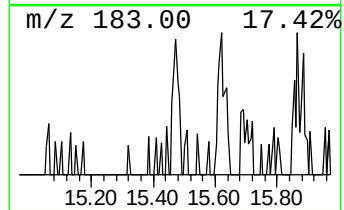
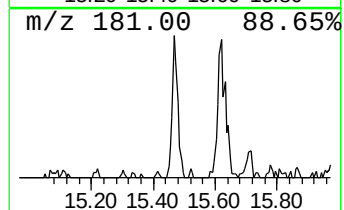
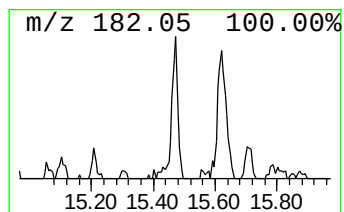
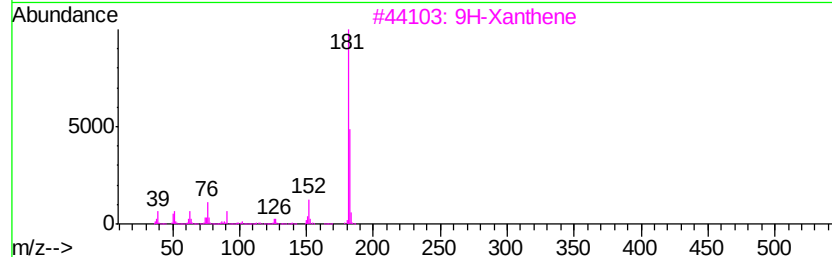
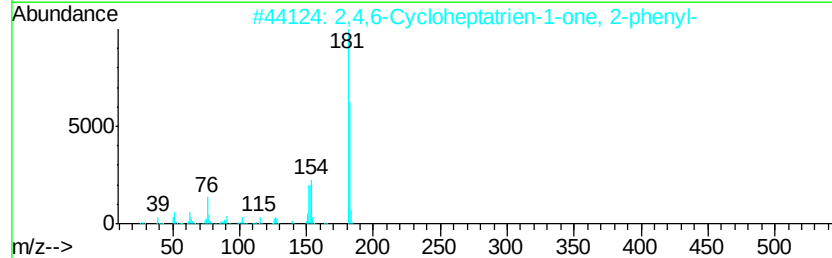
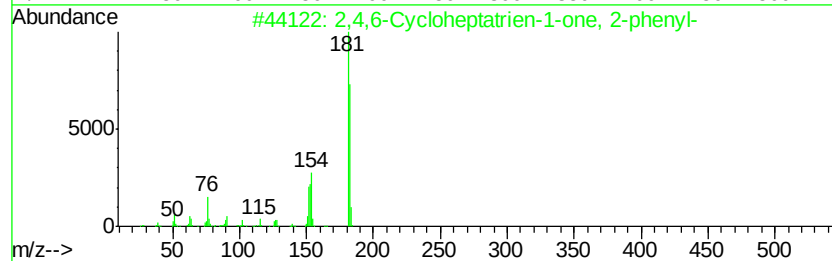
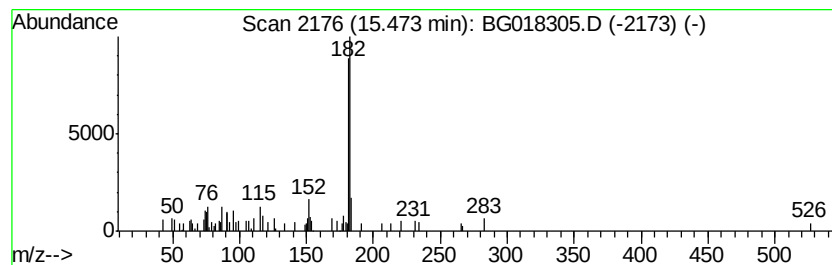
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 2,4,6-Cycloheptatrien-1-one... Concentration Rank 18

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 15.47 | 2.28 ng/ul | 39705 | Acenaphthene-d10 | 14.12 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2,4,6-Cycloheptatrien-1-one, 2-p... | 182 | C13H10O | 014562-09-5 | 74   |
| 2    |    | 2,4,6-Cycloheptatrien-1-one, 2-p... | 182 | C13H10O | 014562-09-5 | 72   |
| 3    |    | 9H-Xanthene                         | 182 | C13H10O | 000092-83-1 | 64   |
| 4    |    | 9H-Fluoren-9-ol                     | 182 | C13H10O | 001689-64-1 | 64   |
| 5    |    | 9H-Xanthene                         | 182 | C13H10O | 000092-83-1 | 64   |



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Quant Title : SVOA CALIBRATION

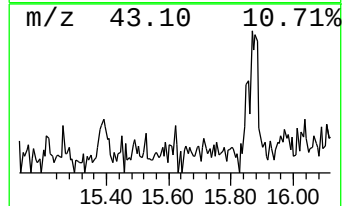
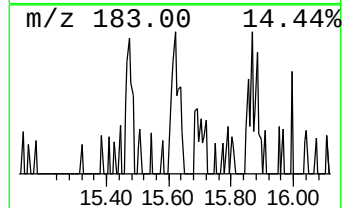
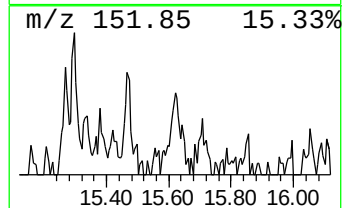
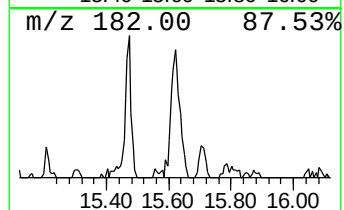
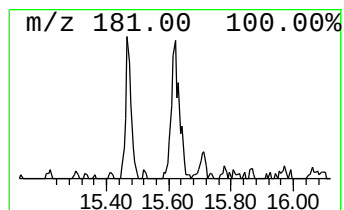
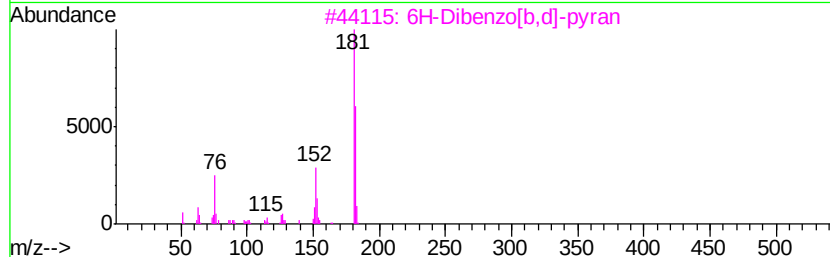
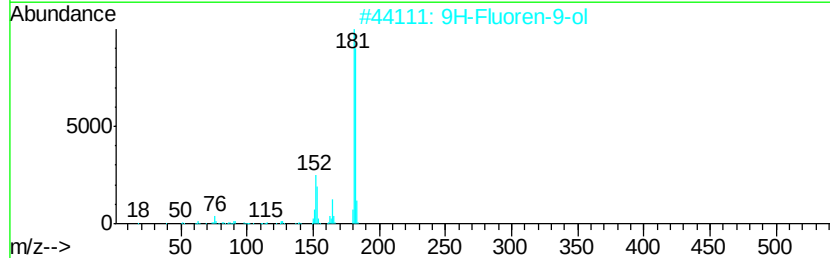
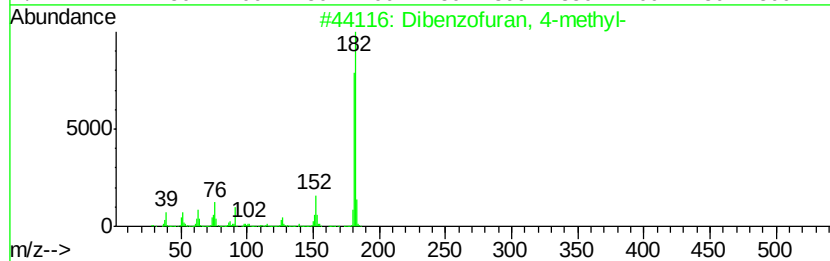
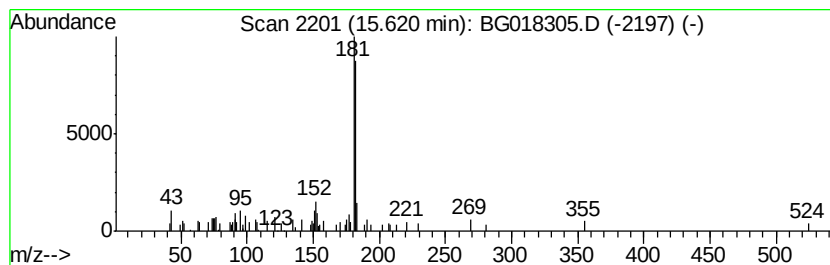
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Dibenzofuran, 4-methyl- Concentration Rank 19

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 15.62 | 2.18 ng/ul | 44509 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Dibenzofuran, 4-methyl-             | 182 | C13H10O  | 007320-53-8 | 87   |
| 2    |    | 9H-Fluoren-9-ol                     | 182 | C13H10O  | 001689-64-1 | 80   |
| 3    |    | 6H-Dibenzo[b,d]-pyran               | 182 | C13H10O  | 000229-95-8 | 72   |
| 4    |    | 9H-Fluoren-9-ol                     | 182 | C13H10O  | 001689-64-1 | 64   |
| 5    |    | Pyridine, 4,4'-(1,2-ethenediyl)bis- | 182 | C12H10N2 | 001135-32-6 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
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Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

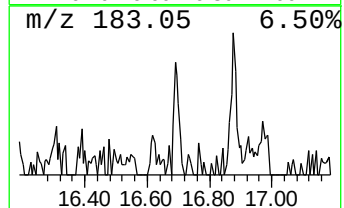
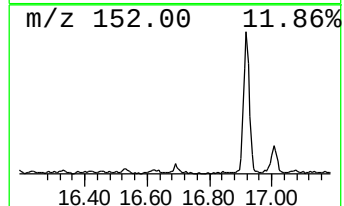
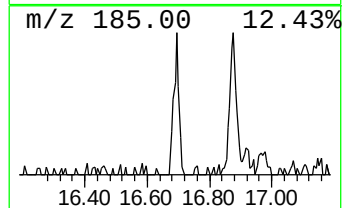
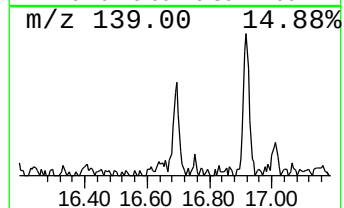
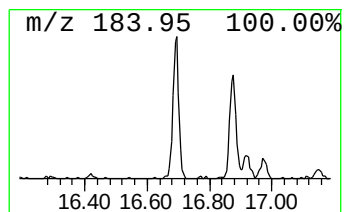
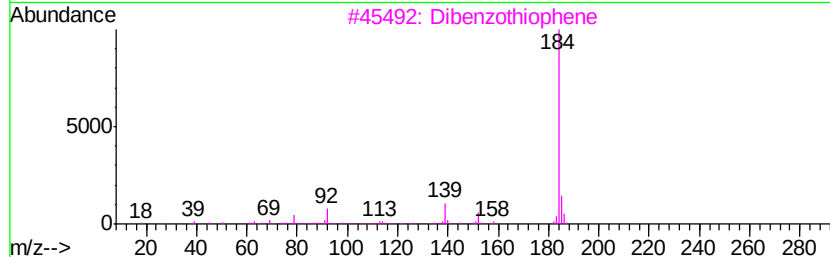
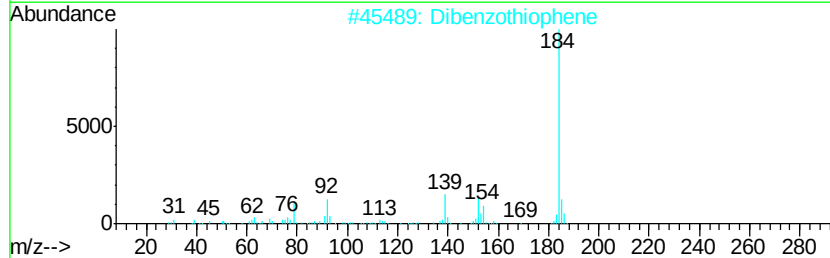
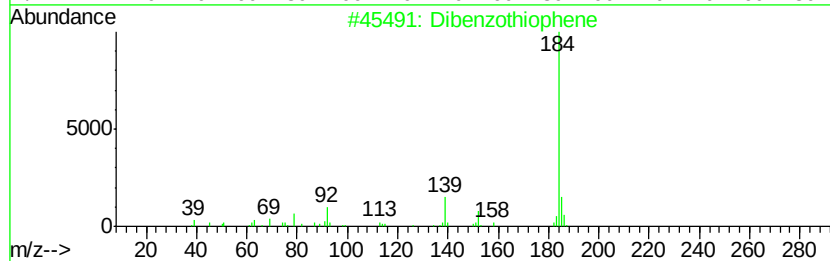
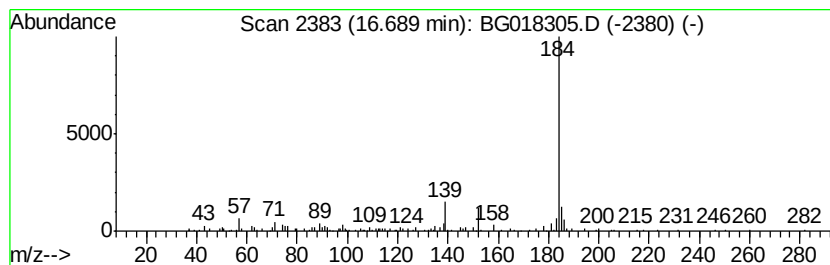
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 Dibenzothiophene Concentration Rank 13

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 16.69 | 3.37 ng/ul | 68756 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Dibenzothiophene        | 184 | C12H8S  | 000132-65-0 | 90   |
| 2    |    | Dibenzothiophene        | 184 | C12H8S  | 000132-65-0 | 87   |
| 3    |    | Dibenzothiophene        | 184 | C12H8S  | 000132-65-0 | 87   |
| 4    |    | Naphtho[2,3-b]thiophene | 184 | C12H8S  | 000268-77-9 | 86   |
| 5    |    | Dibenzothiophene        | 184 | C12H8S  | 000132-65-0 | 83   |





Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleID :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

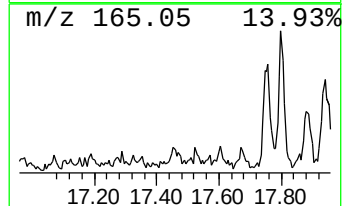
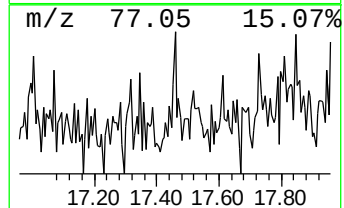
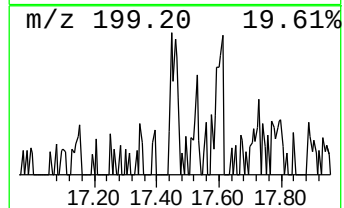
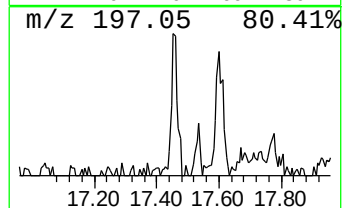
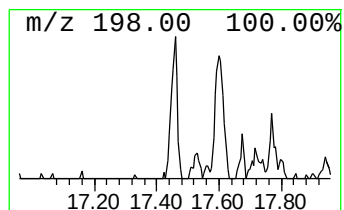
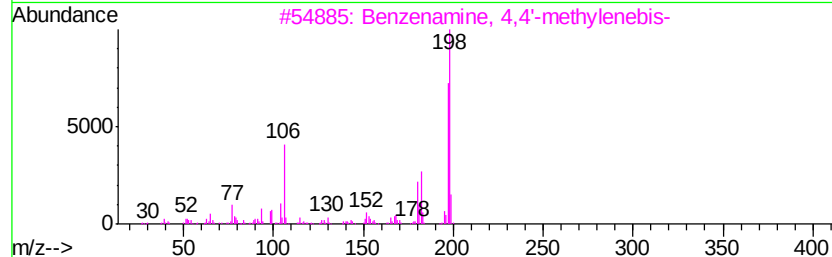
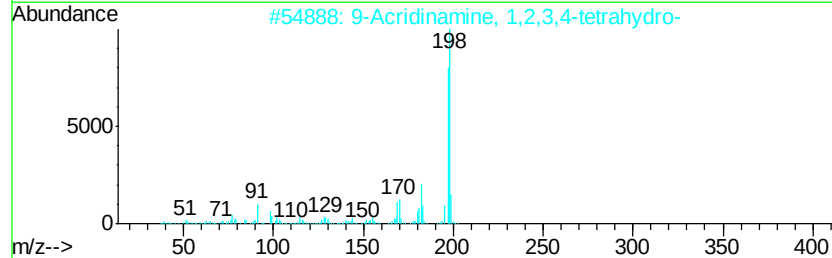
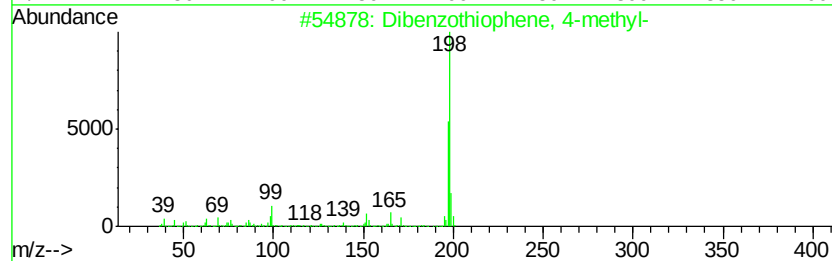
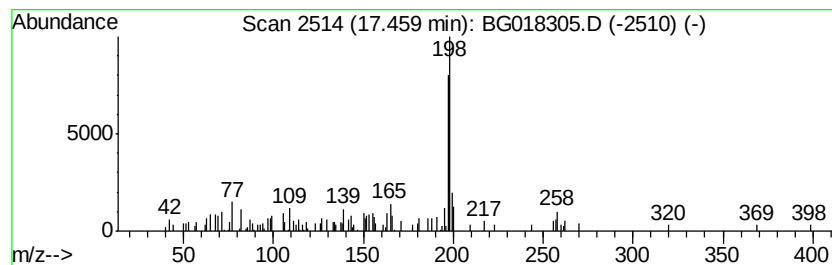
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Dibenzothiophene, 4-methyl- Concentration Rank 17

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 17.46 | 2.34 ng/ul | 47846 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Dibenzothiophene, 4-methyl-         | 198 | C13H10S  | 007372-88-5 | 80   |
| 2    |    | 9-Acridinamine, 1,2,3,4-tetrahydro- | 198 | C13H14N2 | 000321-64-2 | 72   |
| 3    |    | Benzenamine, 4,4'-methylenebis-     | 198 | C13H14N2 | 000101-77-9 | 72   |
| 4    |    | 9-Acridinamine, 1,2,3,4-tetrahydro- | 198 | C13H14N2 | 000321-64-2 | 72   |
| 5    |    | Pyridine, 4-(4-dimethylaminophen... | 198 | C13H14N2 | 001137-80-0 | 64   |





Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

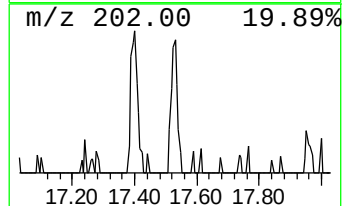
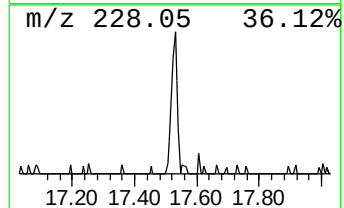
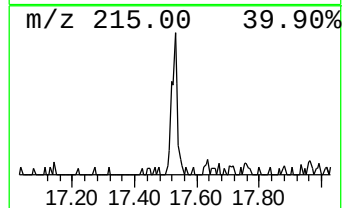
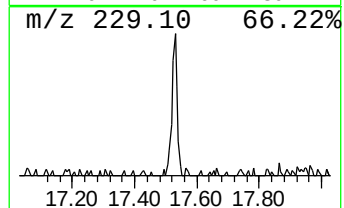
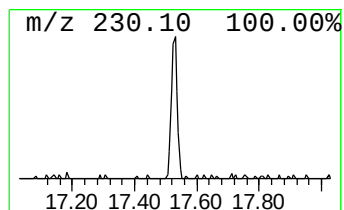
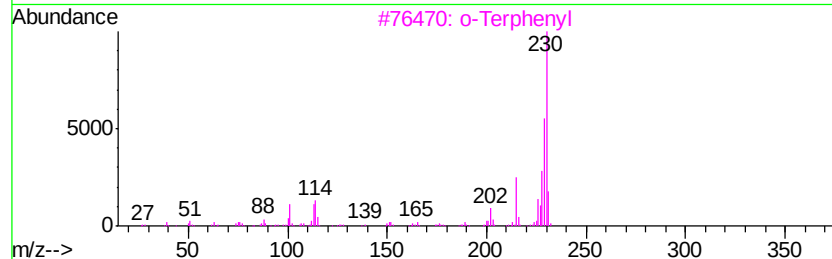
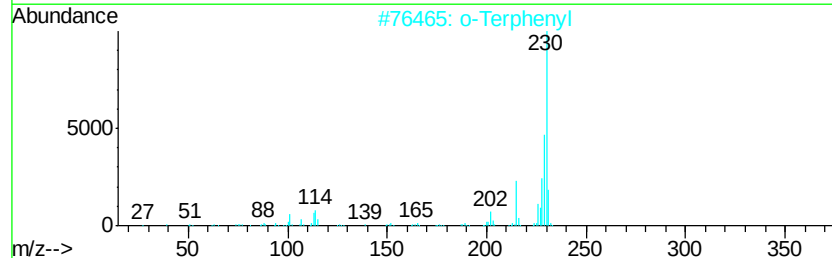
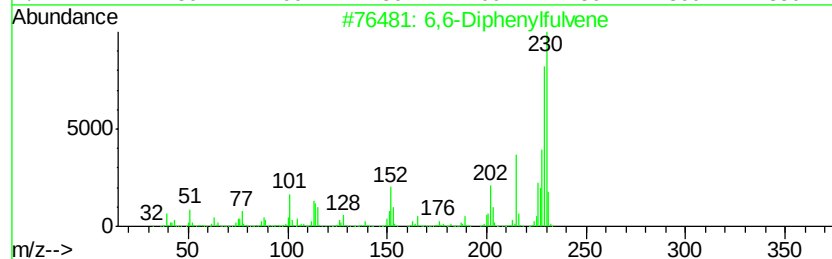
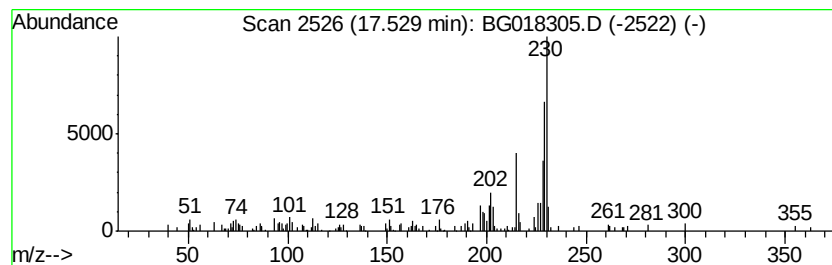
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 6,6-Diphenylfulvene Concentration Rank 16

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 17.53 | 2.67 ng/ul | 54532 | Phenanthrene-d10 | 16.88 |

| Hit# of | 5                                | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|----------------------------------|--------------|-----|---------|-------------|------|
| 1       | 6,6-Diphenylfulvene              |              | 230 | C18H14  | 002175-90-8 | 81   |
| 2       | o-Terphenyl                      |              | 230 | C18H14  | 000084-15-1 | 81   |
| 3       | o-Terphenyl                      |              | 230 | C18H14  | 000084-15-1 | 72   |
| 4       | o-Terphenyl                      |              | 230 | C18H14  | 000084-15-1 | 68   |
| 5       | Benz[a]anthracene, 7,12-dihydro- |              | 230 | C18H14  | 016434-59-6 | 58   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

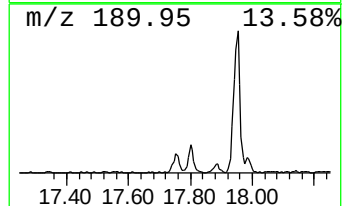
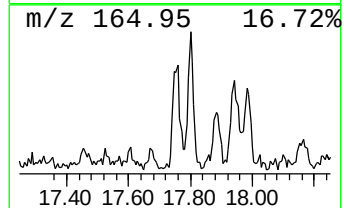
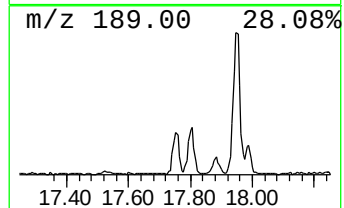
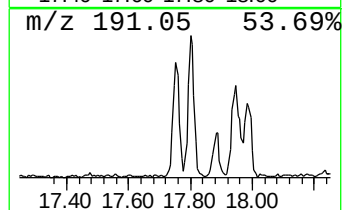
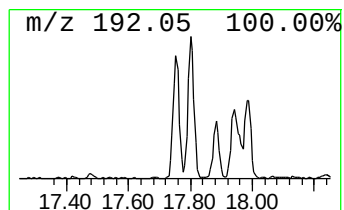
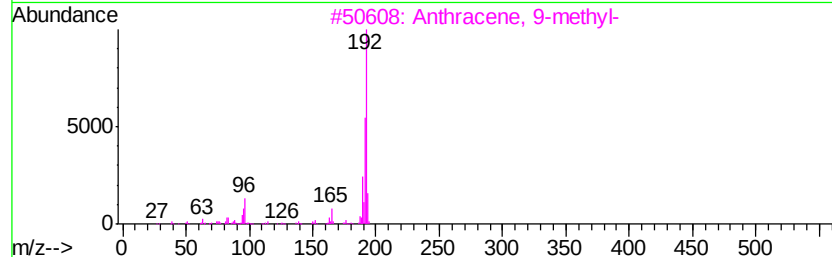
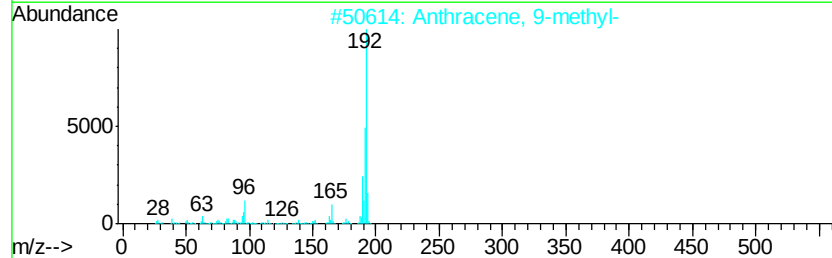
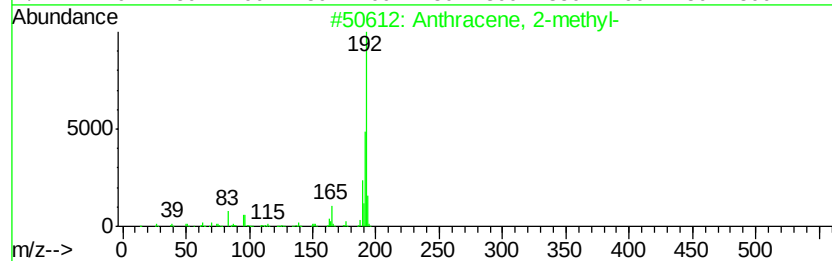
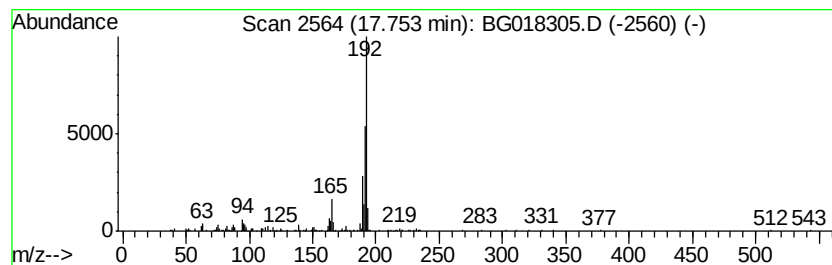
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Anthracene, 2-methyl- Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.75 | 6.13 ng/ul | 125300 | Phenanthrene-d10 | 16.88 |

| Hit# of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------|-----|---------|-------------|------|
| 1       |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 90   |
| 2       |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 81   |
| 3       |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 81   |
| 4       |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 81   |
| 5       |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 76   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

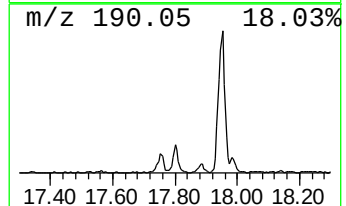
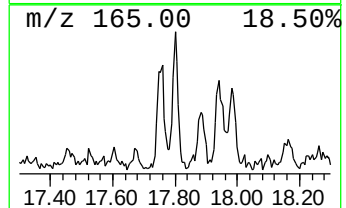
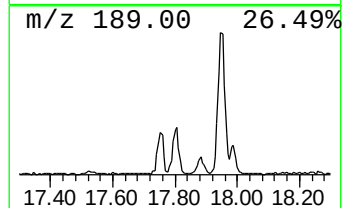
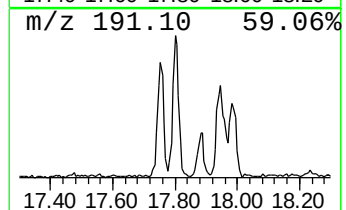
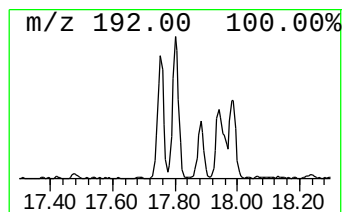
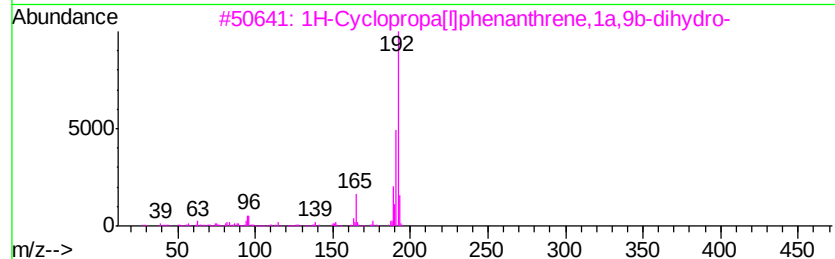
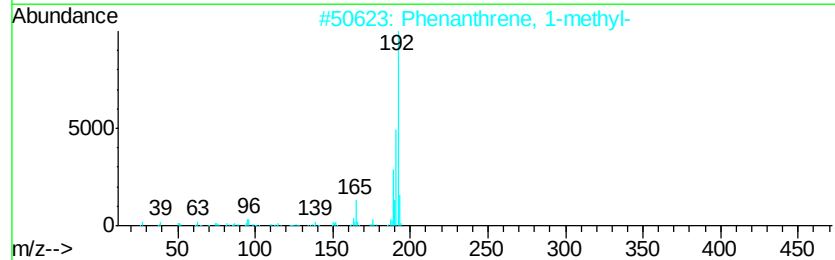
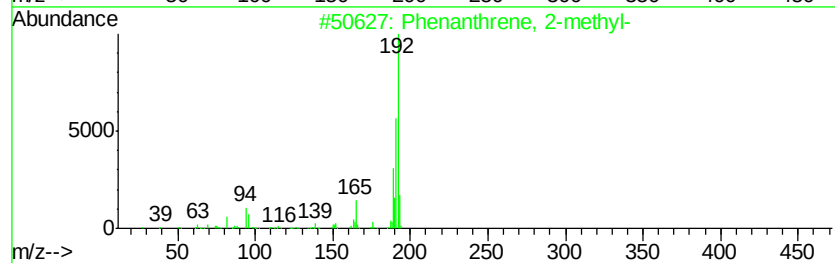
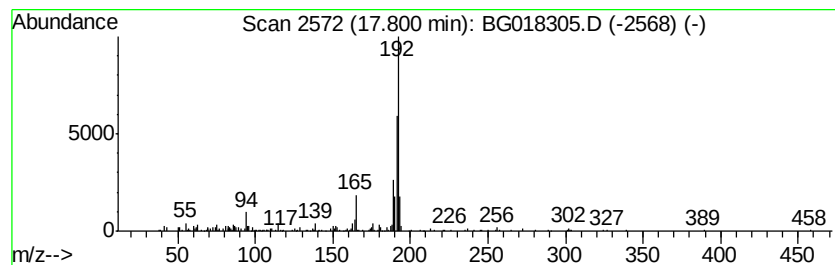
Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 4

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 17.80     | 9.58 ng/ul                          | 195768 | Phenanthrene-d10 | 16.88       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 91   |
| 2         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 90   |
| 3         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 90   |
| 4         | Naphtho[2,3-b]norborene             | 192    | C15H12           | 107426-38-0 | 89   |
| 5         | Anthracene, 2-methyl-               | 192    | C15H12           | 000613-12-7 | 87   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

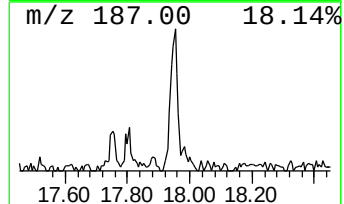
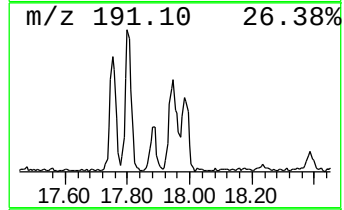
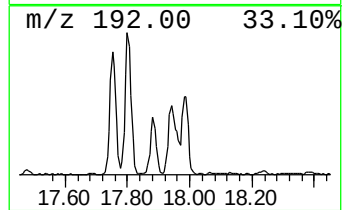
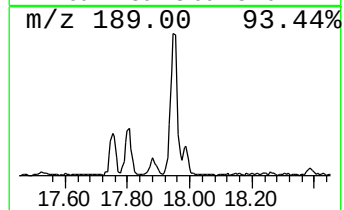
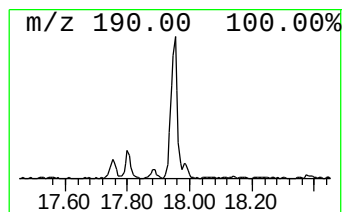
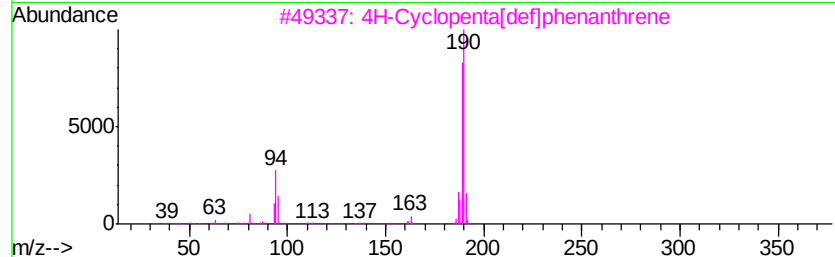
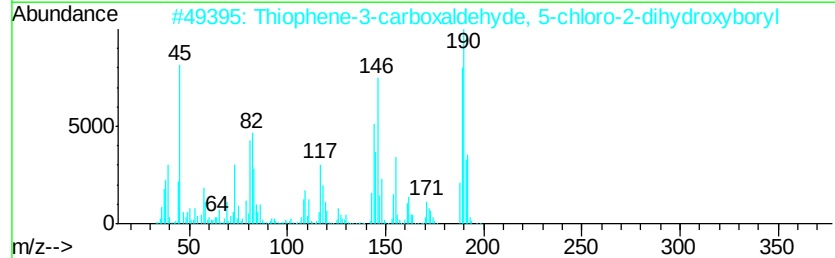
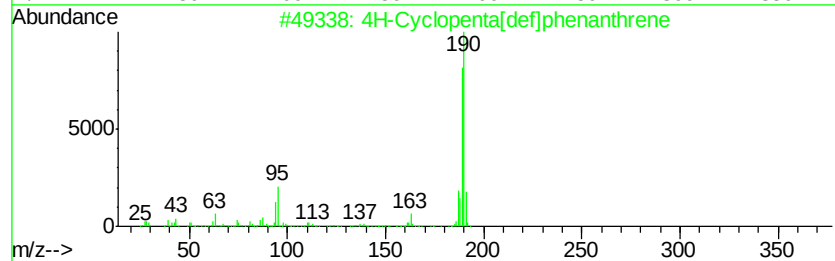
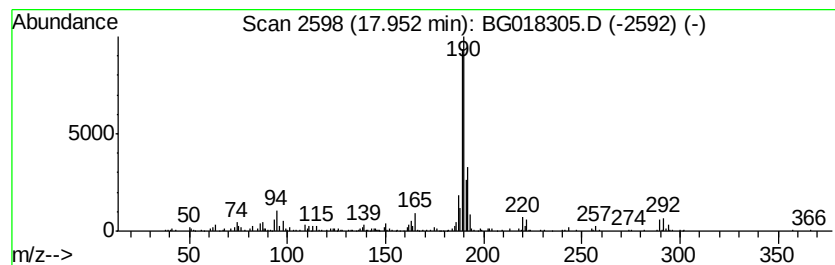
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 17.95 | 14.24 ng/ul | 291032 | Phenanthrene-d10 | 16.88 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|---------|---|-------------------------------------|-----|------------|-------------|------|
| 1       |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10     | 000203-64-5 | 64   |
| 2       |   | Thiophene-3-carboxaldehyde, 5-ch... | 190 | C5H4BClO3S | 036155-87-0 | 64   |
| 3       |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10     | 000203-64-5 | 50   |
| 4       |   | 6H-Cyclobuta[flk]phenanthrene       | 190 | C15H10     | 083469-43-6 | 50   |
| 5       |   | 2,3,5,6-Tetramethylterephthald...   | 190 | C12H14O2   | 007072-01-7 | 38   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

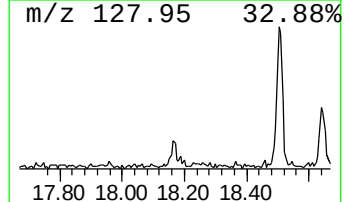
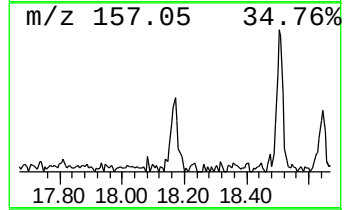
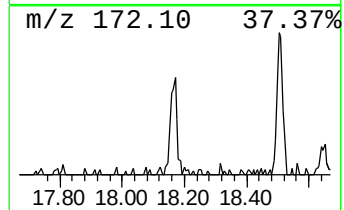
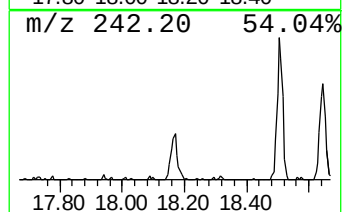
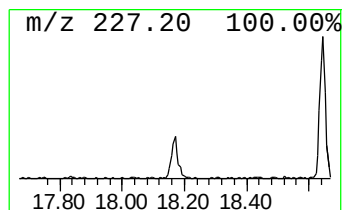
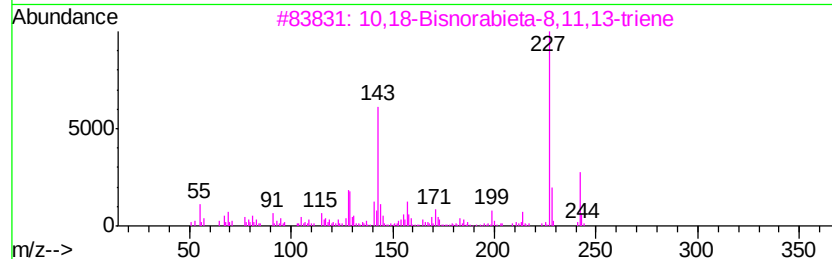
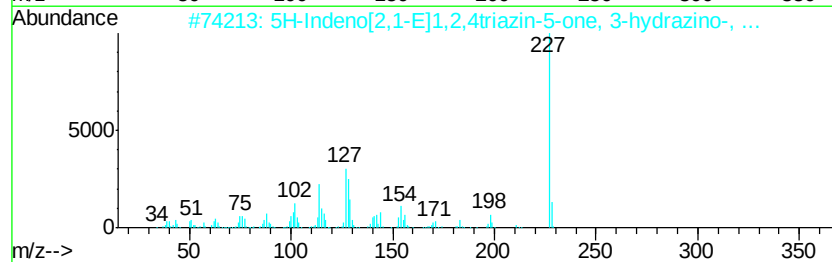
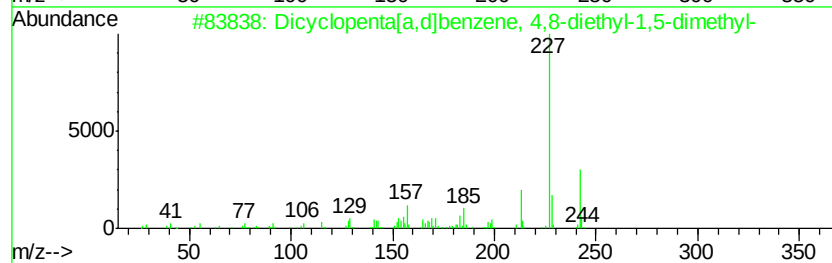
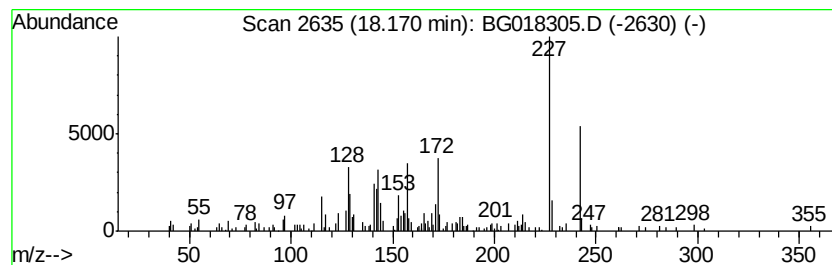
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Dicyclopenta[a,d]benzene, 4... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.17 | 5.04 ng/ul | 102976 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Dicyclopenta[a,d]benzene, 4,8-di... | 242 | C18H26  | 1000156-41-6 | 50   |
| 2    |    | 5H-Indeno[2,1-E]1,2,4triazin-5-o... | 227 | C10H9N7 | 1000271-75-1 | 50   |
| 3    |    | 10,18-Bisnorabieta-8,11,13-triene   | 242 | C18H26  | 032624-67-2  | 43   |
| 4    |    | Benzocyclopentene, 4,6,7-triethy... | 242 | C18H26  | 1000156-41-5 | 41   |
| 5    |    | S-Indacene, 1,2,3,5,6,7-hexahydr... | 242 | C18H26  | 017465-58-6  | 27   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

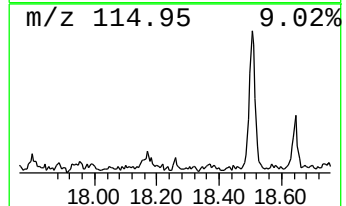
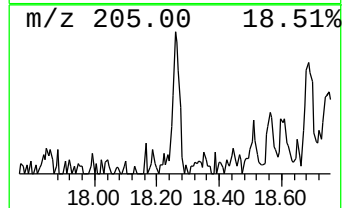
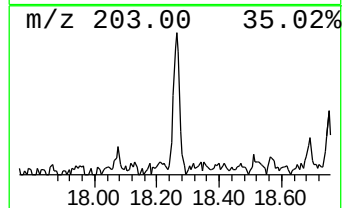
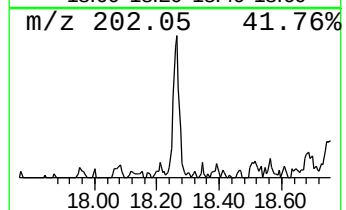
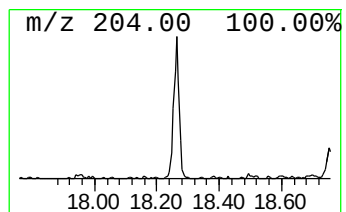
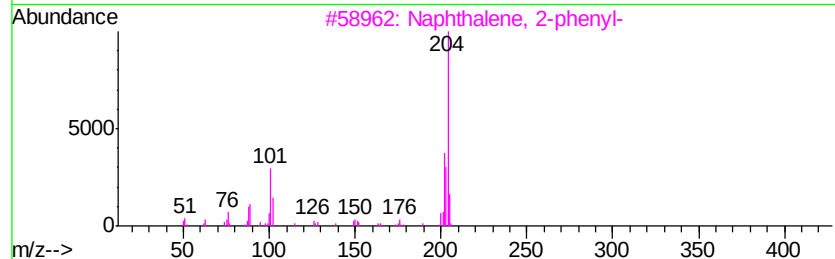
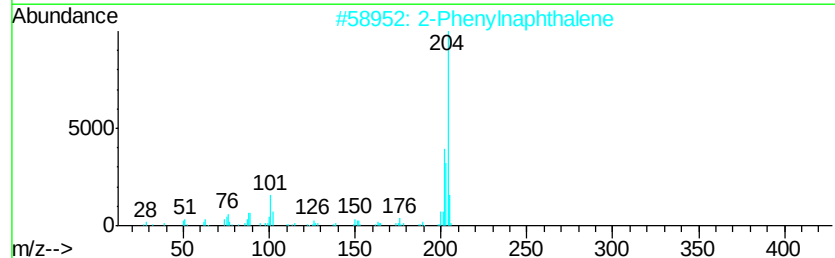
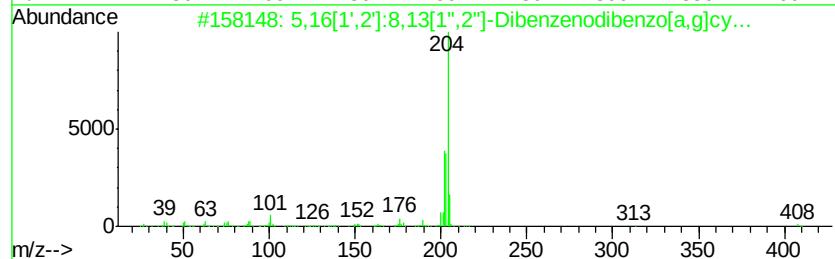
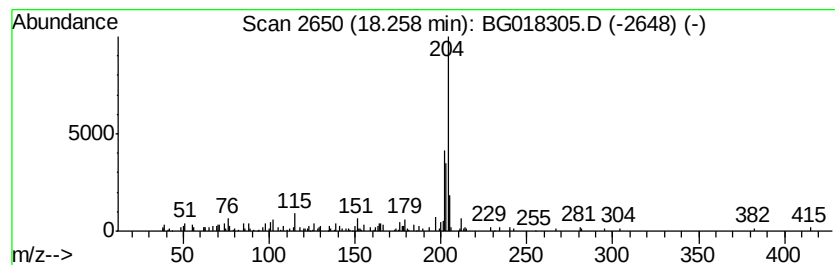
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 15

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 18.26 | 3.11 ng/ul | 63537 | Phenanthrene-d10 | 16.88 |

| Hit# | of                                  | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24       | 005672-97-9 | 87      |      |      |
| 2    | 2-Phenylnaphthalene                 | 204 | C16H12       | 035465-71-5 | 76      |      |      |
| 3    | Naphthalene, 2-phenyl-              | 204 | C16H12       | 000612-94-2 | 58      |      |      |
| 4    | Naphthalene, 2-phenyl-              | 204 | C16H12       | 000612-94-2 | 58      |      |      |
| 5    | 6-Cyanophenanthridine               | 204 | C14H8N2      | 002176-28-5 | 58      |      |      |







Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

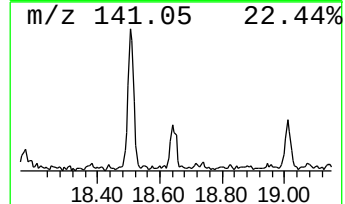
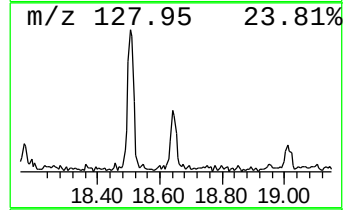
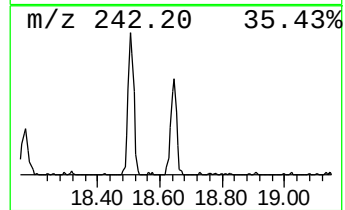
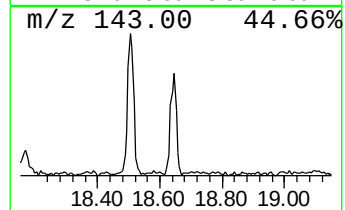
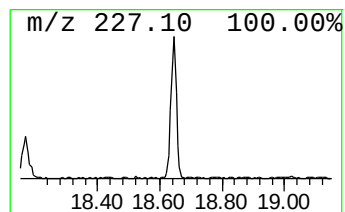
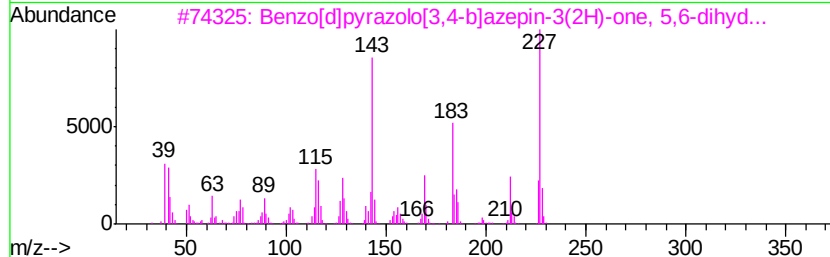
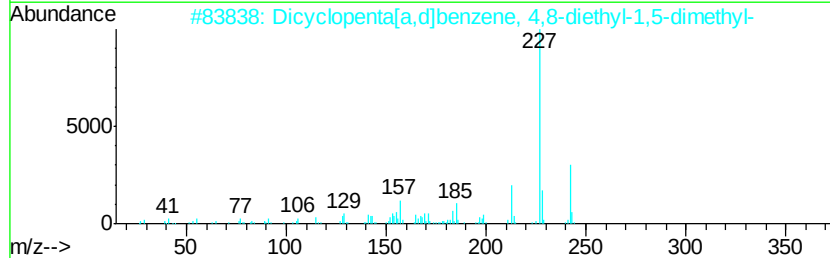
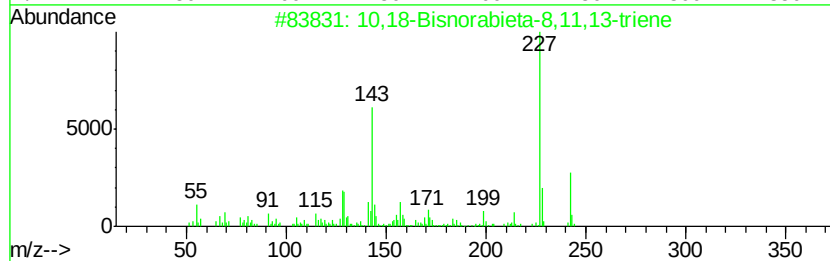
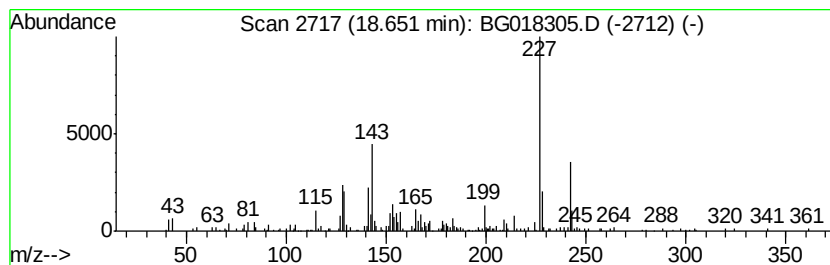
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 10,18-Bisnorabieta-8,11,13-... Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.65 | 8.51 ng/ul | 173896 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 10,18-Bisnorabieta-8,11,13-triene   | 242 | C18H26     | 032624-67-2  | 93   |
| 2    |    | Dicyclopenta[a,d]benzene, 4,8-di... | 242 | C18H26     | 1000156-41-6 | 83   |
| 3    |    | Benzo[d]pyrazolo[3,4-b]azepin-3(... | 227 | C13H13N3O  | 263550-89-6  | 70   |
| 4    |    | Styrene, 2,3,5,6-tetraethyl-4-vi... | 242 | C18H26     | 1000156-41-3 | 50   |
| 5    |    | 7-Diethylamino-2-oxo-2H-chromene... | 242 | C14H14N2O2 | 332411-59-3  | 49   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

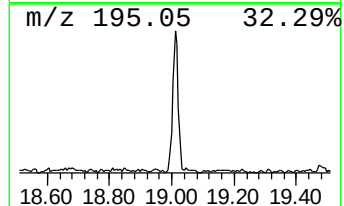
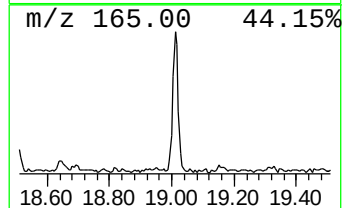
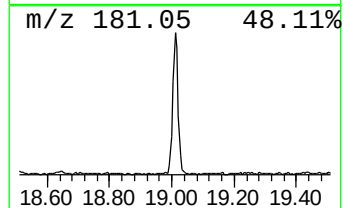
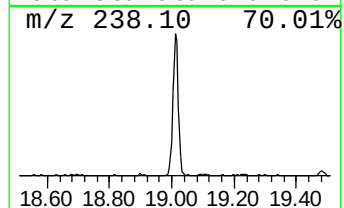
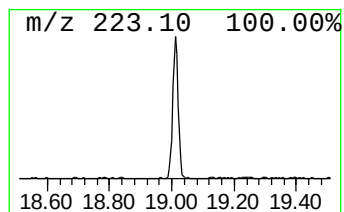
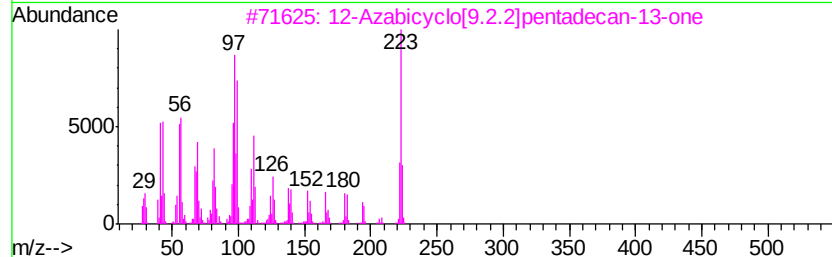
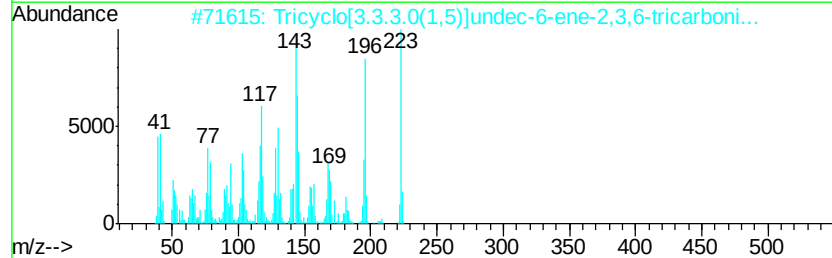
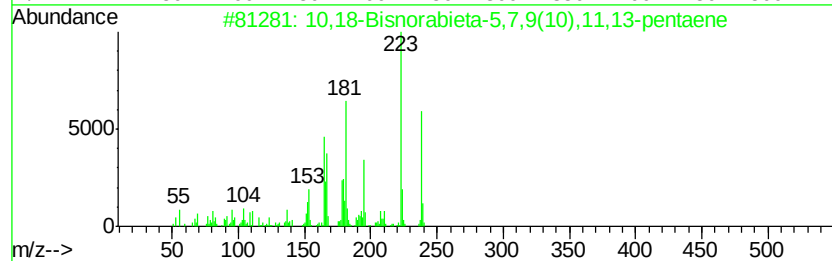
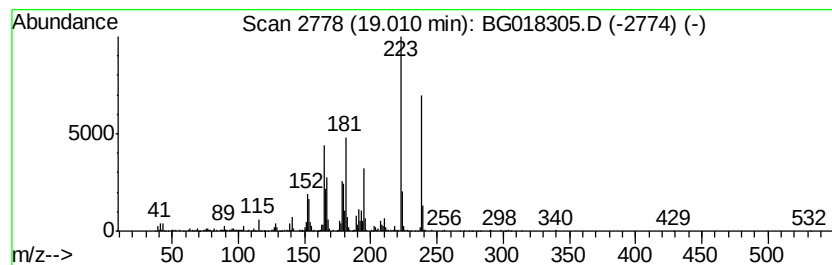
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 10,18-Bisnorabieta-5,7,9(10)... Concentration Rank 2

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 19.01 | 23.71 ng/ul | 617145 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 10,18-Bisnorabieta-5,7,9(10),11,... | 238 | C18H22   | 006566-19-4  | 99   |
| 2    |    | Tricyclo[3.3.3.0(1,5)]undec-6-en... | 223 | C14H13N3 | 118894-71-6  | 64   |
| 3    |    | 12-Azabicyclo[9.2.2]pentadecan-1... | 223 | C14H25NO | 1000191-26-7 | 60   |
| 4    |    | 4,4'-Diisopropylbiphenyl            | 238 | C18H22   | 018970-30-4  | 55   |
| 5    |    | 4,4'-Diisopropylbiphenyl            | 238 | C18H22   | 018970-30-4  | 50   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

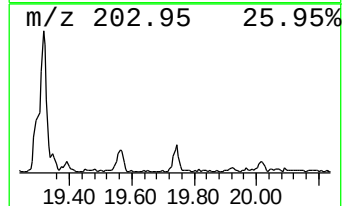
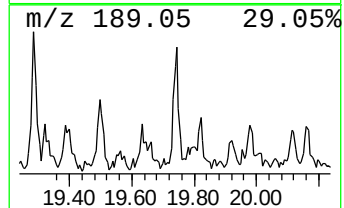
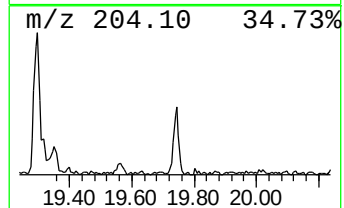
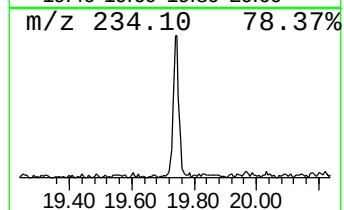
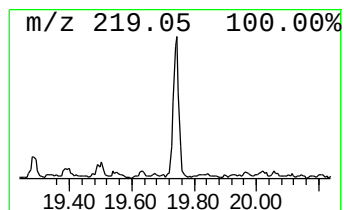
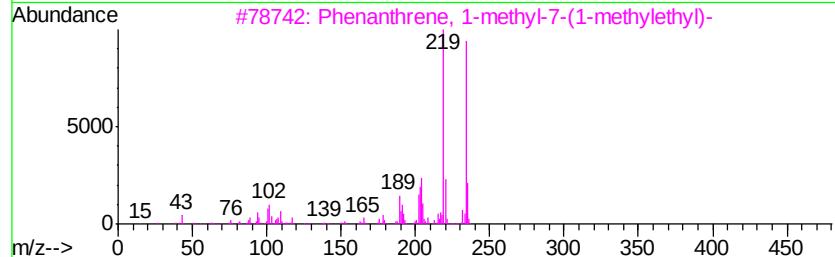
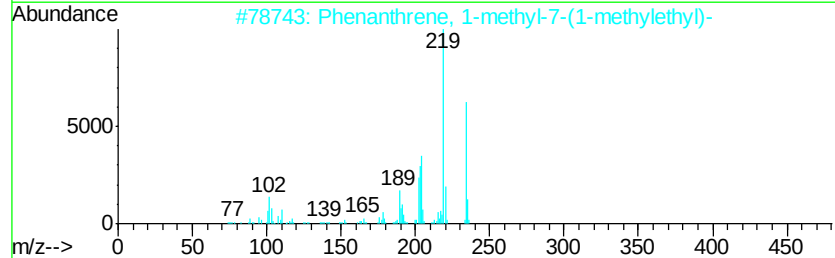
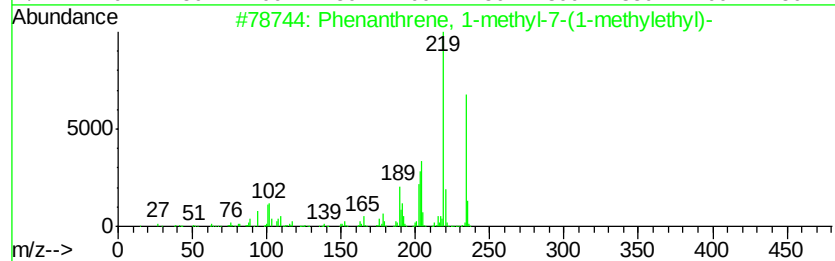
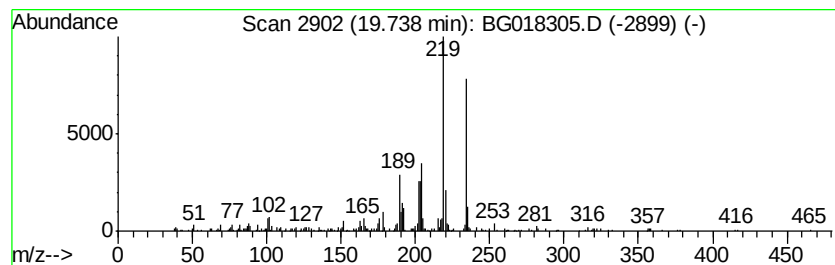
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Phenanthrene, 1-methyl-7-(1... Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.74 | 5.80 ng/ul | 150883 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 1-methyl-7-(1-meth... | 234 | C18H18  | 000483-65-8 | 95   |
| 2    |    | Phenanthrene, 1-methyl-7-(1-meth... | 234 | C18H18  | 000483-65-8 | 95   |
| 3    |    | Phenanthrene, 1-methyl-7-(1-meth... | 234 | C18H18  | 000483-65-8 | 94   |
| 4    |    | 2-Isopropyl-10-methylphenanthrene   | 234 | C18H18  | 066552-97-4 | 93   |
| 5    |    | Phenanthrene, 2,4,5,7-tetramethyl-  | 234 | C18H18  | 007396-38-5 | 93   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

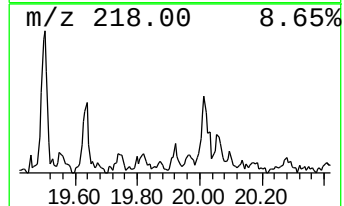
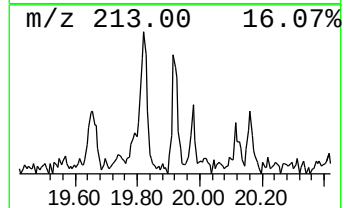
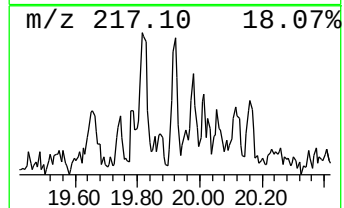
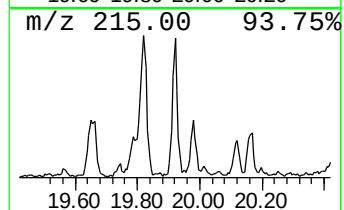
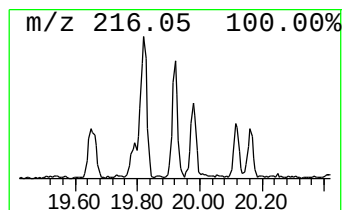
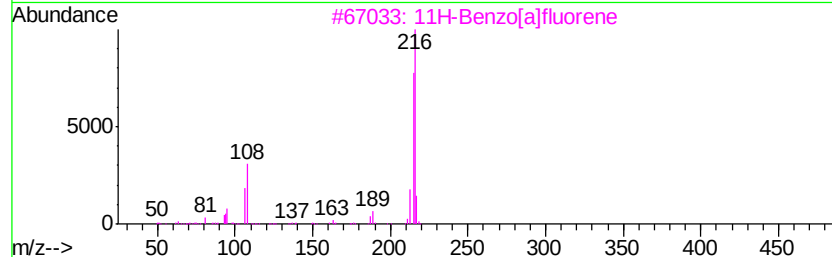
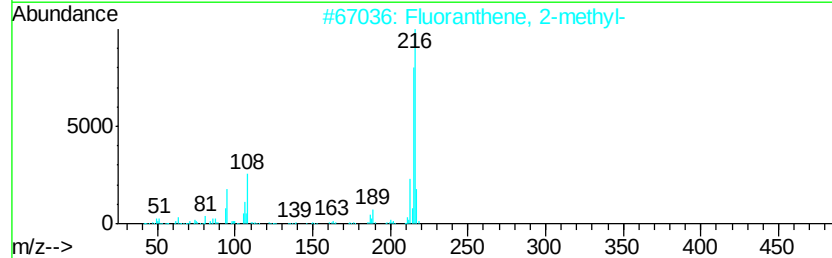
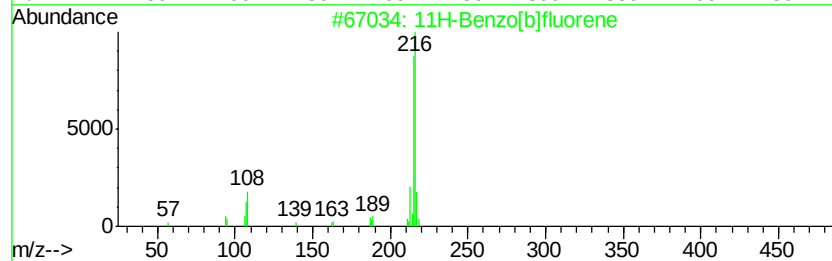
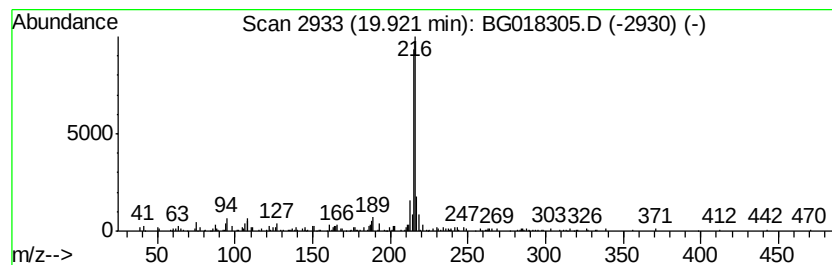
Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 14

| R.T.    | EstConc    | Area                    | Relative to ISTD | R.T.           |
|---------|------------|-------------------------|------------------|----------------|
| 19.92   | 3.20 ng/ul | 83189                   | Chrysene-d12     | 21.09          |
| Hit# of | 5          | Tentative ID            | MW MolForm       | CAS# Qual      |
| 1       |            | 11H-Benzo[b]fluorene    | 216 C17H12       | 000243-17-4 93 |
| 2       |            | Fluoranthene, 2-methyl- | 216 C17H12       | 033543-31-6 91 |
| 3       |            | 11H-Benzo[a]fluorene    | 216 C17H12       | 000238-84-6 90 |
| 4       |            | 11H-Benzo[b]fluorene    | 216 C17H12       | 000243-17-4 90 |
| 5       |            | Pyrene, 1-methyl-       | 216 C17H12       | 002381-21-7 90 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

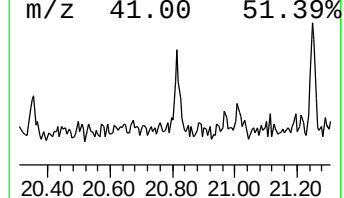
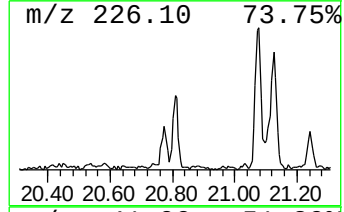
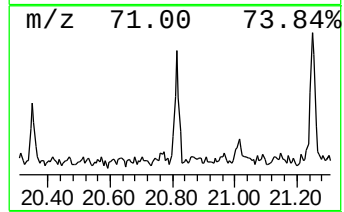
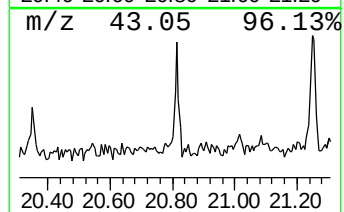
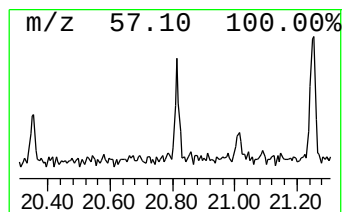
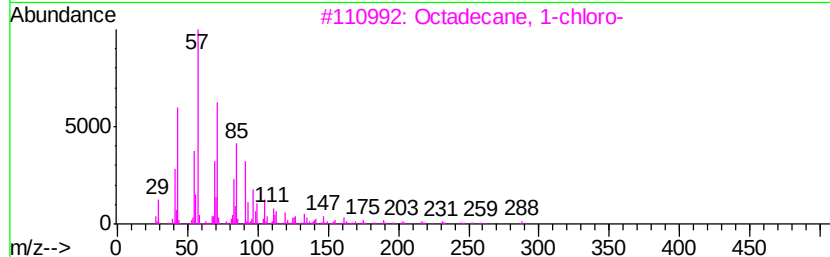
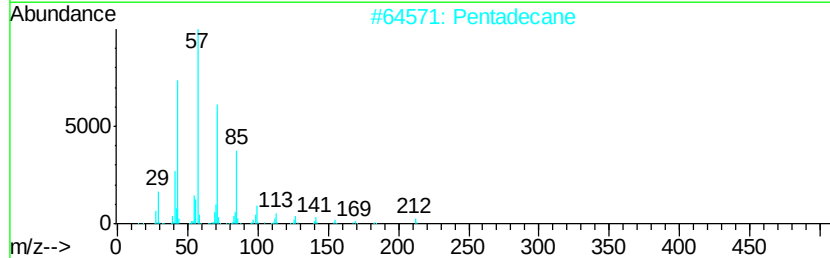
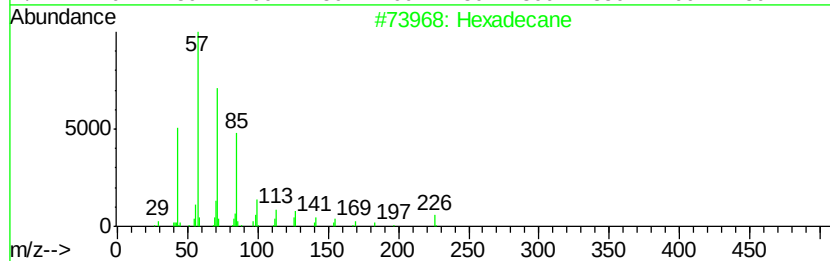
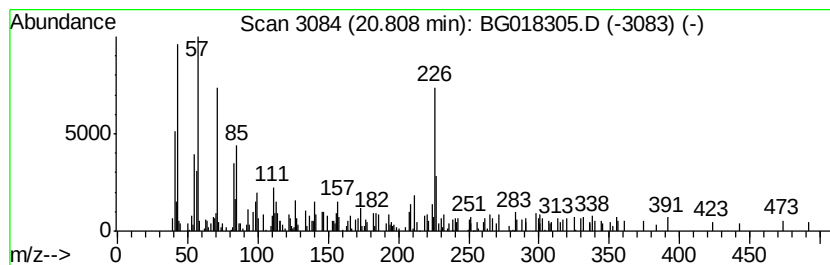
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.81 | 4.03 ng/ul | 105004 | Chrysene-d12     | 21.09 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------|-----|----------|-------------|------|
| 1    |    |   | Hexadecane             | 226 | C16H34   | 000544-76-3 | 74   |
| 2    |    |   | Pentadecane            | 212 | C15H32   | 000629-62-9 | 70   |
| 3    |    |   | Octadecane, 1-chloro-  | 288 | C18H37Cl | 003386-33-2 | 70   |
| 4    |    |   | Eicosane               | 282 | C20H42   | 000112-95-8 | 64   |
| 5    |    |   | Heptadecane, 4-methyl- | 254 | C18H38   | 026429-11-8 | 64   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

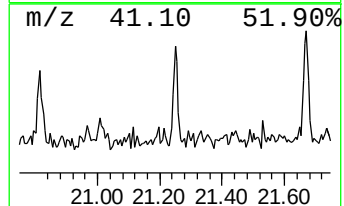
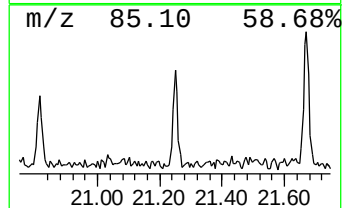
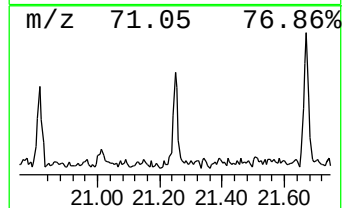
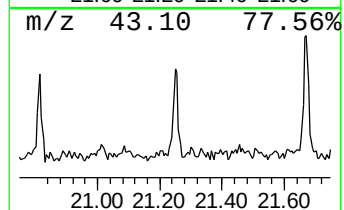
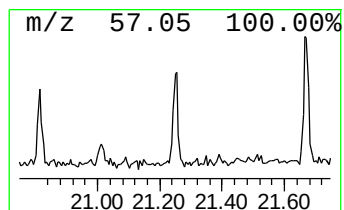
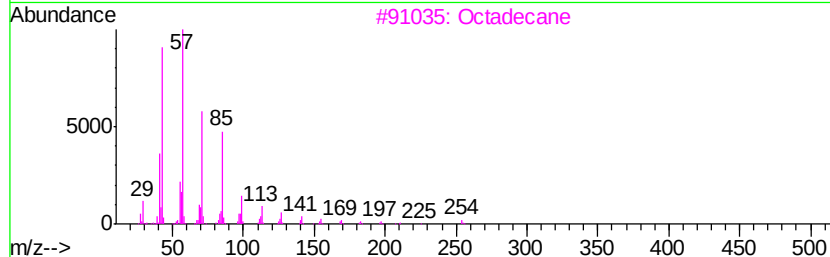
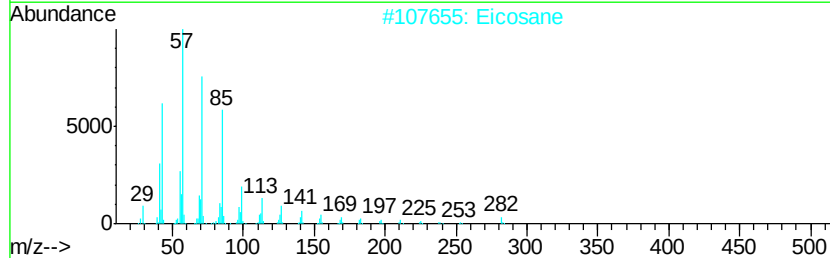
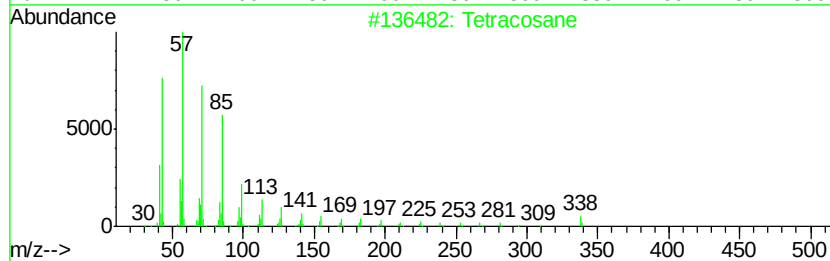
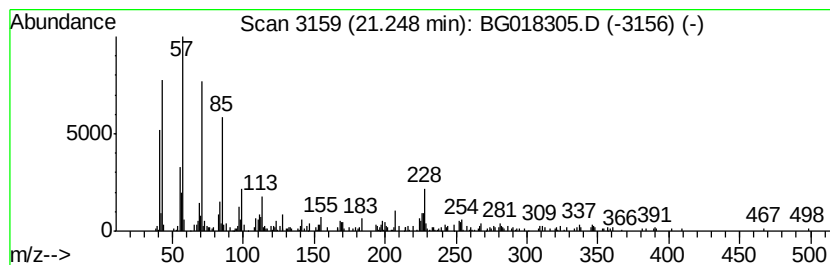
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.25 | 4.76 ng/ul | 123807 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Tetracosane  | 338 | C24H50  | 000646-31-1 | 93   |
| 2    |    | Eicosane     | 282 | C20H42  | 000112-95-8 | 92   |
| 3    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 89   |
| 4    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 89   |
| 5    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 89   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

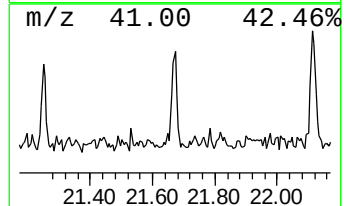
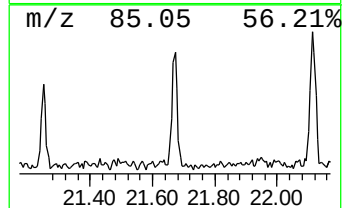
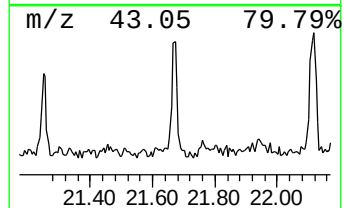
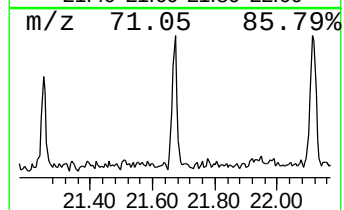
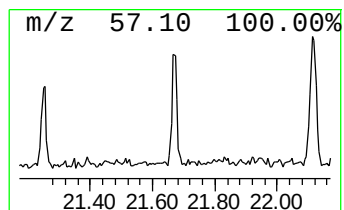
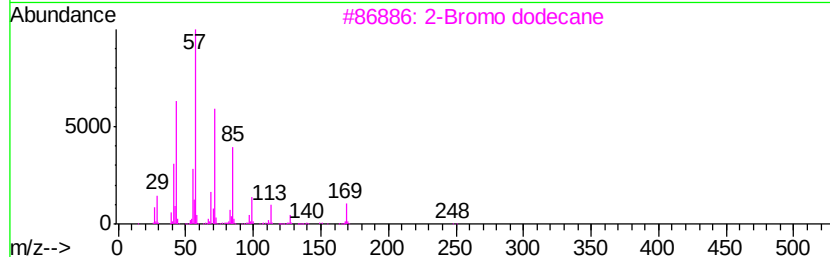
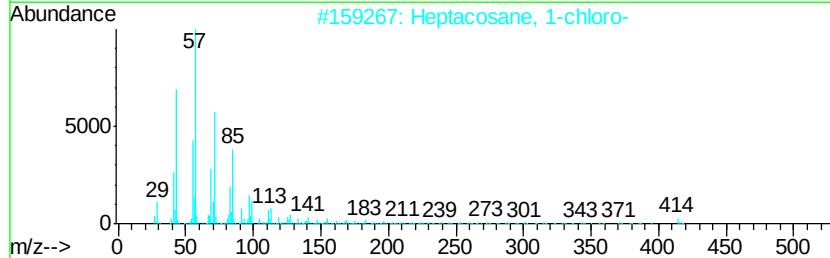
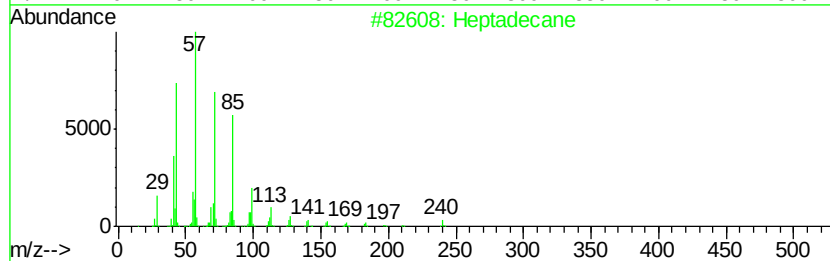
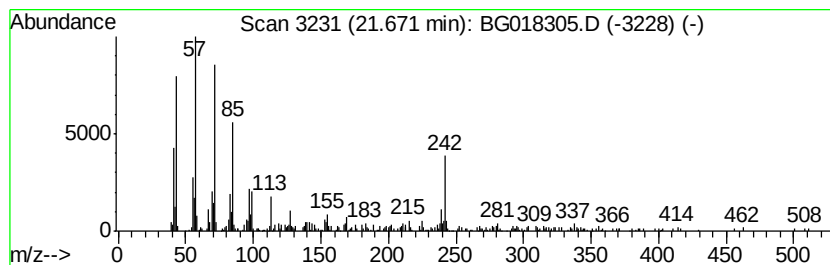
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.67 | 6.46 ng/ul | 168091 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------|-----|----------|-------------|------|
| 1    | 5  | Heptadecane            | 240 | C17H36   | 000629-78-7 | 91   |
| 2    |    | Heptacosane, 1-chloro- | 414 | C27H55Cl | 062016-79-9 | 89   |
| 3    |    | 2-Bromo dodecane       | 248 | C12H25Br | 013187-99-0 | 83   |
| 4    |    | Nonadecane, 2-methyl-  | 282 | C20H42   | 001560-86-7 | 83   |
| 5    |    | Heptacosane            | 380 | C27H56   | 000593-49-7 | 83   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080615\  
Data File : BG018305.D  
Acq On : 6 Aug 2015 15:20  
Operator : UM/TP  
Sample : G3109-15DL 5X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W3DL

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

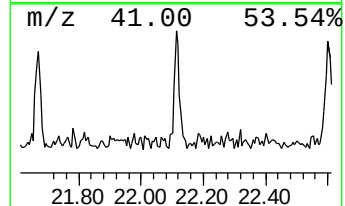
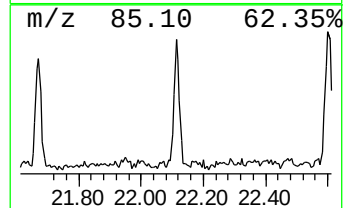
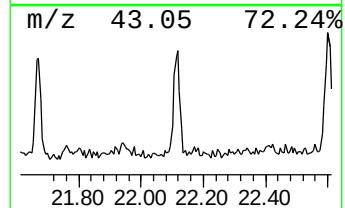
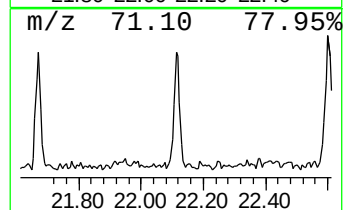
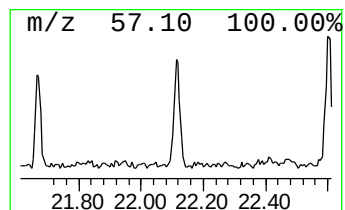
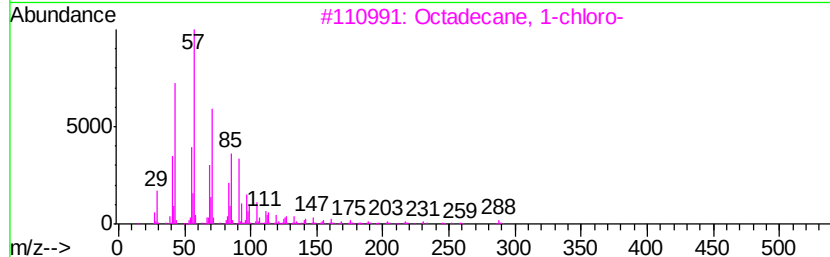
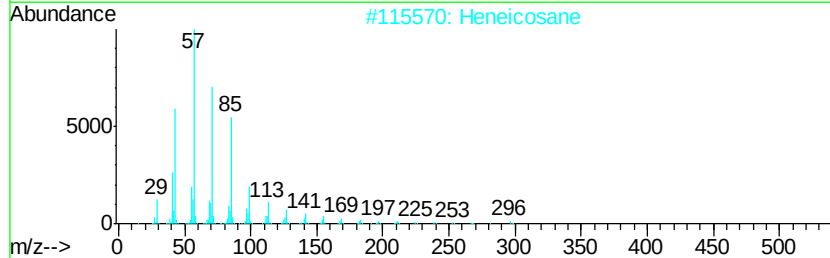
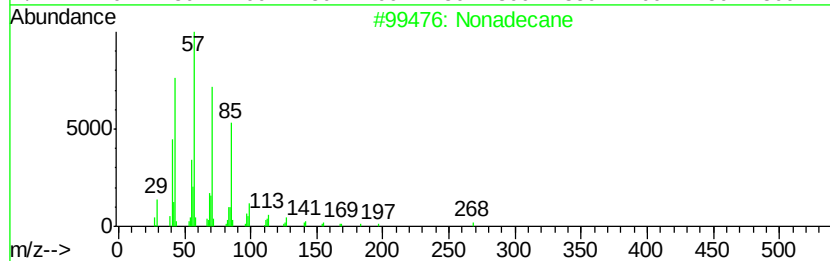
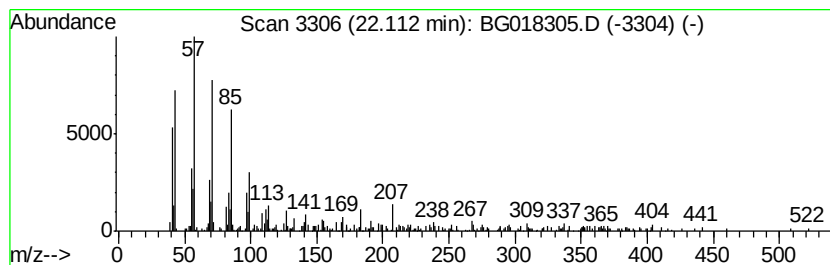
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.11 | 7.89 ng/ul | 205464 | Chrysene-d12     | 21.09 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-----------------------|-----|----------|-------------|------|
| 1    |    |   | Nonadecane            | 268 | C19H40   | 000629-92-5 | 94   |
| 2    |    |   | Heneicosane           | 296 | C21H44   | 000629-94-7 | 94   |
| 3    |    |   | Octadecane, 1-chloro- | 288 | C18H37Cl | 003386-33-2 | 89   |
| 4    |    |   | Heptacosane           | 380 | C27H56   | 000593-49-7 | 87   |
| 5    |    |   | Octadecane, 1-chloro- | 288 | C18H37Cl | 003386-33-2 | 87   |





Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG080615\  
 Data File : BG018305.D  
 Acq On : 6 Aug 2015 15:20  
 Operator : UM/TP  
 Sample : G3109-15DL 5X  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C01W3DL

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG080515.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 |
| 2,4,6-Cycloheptat... | 15.47 | 2.3     | ng/ul | 39705    | 3                     | 14.12 | 348759 | 20.0 |
| Dibenzofuran, 4-m... | 15.62 | 2.2     | ng/ul | 44509    | 4                     | 16.88 | 408629 | 20.0 |
| Dibenzothiophene     | 16.69 | 3.4     | ng/ul | 68756    | 4                     | 16.88 | 408629 | 20.0 |
| Dibenzothiophene,... | 17.46 | 2.3     | ng/ul | 47846    | 4                     | 16.88 | 408629 | 20.0 |
| 6,6-Diphenylfulvene  | 17.53 | 2.7     | ng/ul | 54532    | 4                     | 16.88 | 408629 | 20.0 |
| Anthracene, 2-met... | 17.75 | 6.1     | ng/ul | 125300   | 4                     | 16.88 | 408629 | 20.0 |
| Phenanthrene, 2-m... | 17.80 | 9.6     | ng/ul | 195768   | 4                     | 16.88 | 408629 | 20.0 |
| 4H-Cyclopenta[def... | 17.95 | 14.2    | ng/ul | 291032   | 4                     | 16.88 | 408629 | 20.0 |
| Dicyclopenta[a,d]... | 18.17 | 5.0     | ng/ul | 102976   | 4                     | 16.88 | 408629 | 20.0 |
| 5,16[1',2']:8,13[... | 18.26 | 3.1     | ng/ul | 63537    | 4                     | 16.88 | 408629 | 20.0 |
| 4b,8-Dimethyl-2-i... | 18.50 | 36.7    | ng/ul | 749306   | 4                     | 16.88 | 408629 | 20.0 |
| 10,18-Bisnorabiet... | 18.65 | 8.5     | ng/ul | 173896   | 4                     | 16.88 | 408629 | 20.0 |
| 10,18-Bisnorabiet... | 19.01 | 23.7    | ng/ul | 617145   | 5                     | 21.09 | 520607 | 20.0 |
| Phenanthrene, 1-m... | 19.74 | 5.8     | ng/ul | 150883   | 5                     | 21.09 | 520607 | 20.0 |
| 11H-Benzo[b]fluorene | 19.92 | 3.2     | ng/ul | 83189    | 5                     | 21.09 | 520607 | 20.0 |
| (DEL) Alkane: Str... | 20.81 | 4.0     | ng/ul | 105004   | 5                     | 21.09 | 520607 | 20.0 |
| (DEL) Alkane: Str... | 21.25 | 4.8     | ng/ul | 123807   | 5                     | 21.09 | 520607 | 20.0 |
| (DEL) Alkane: Str... | 21.67 | 6.5     | ng/ul | 168091   | 5                     | 21.09 | 520607 | 20.0 |
| (DEL) Alkane: Str... | 22.11 | 7.9     | ng/ul | 205464   | 5                     | 21.09 | 520607 | 20.0 |